Component RT	CAS#	Compound Name	Formul	a
4.3249	67-71-0	Dimethyl sulfone	C2H6C)2S.
4.5005	10076-48-9	Methyl pyruvate dimethyl acetal	C6H12	204
4.5773	927-83-3	Diazene, bis(1,1-dimethylethyl)-	C8H18	
4.6390	136-36-7	1,3-Benzenediol, monobenzoate	C13H	
4.7784	624-45-3	Pentanoic acid, 4-oxo-, methyl ester	C6H1	
4.9599	927-83-3	Diazene, bis(1,1-dimethylethyl)-	C8H1	
5.0248	16747-26-5	Hexane, 2,2,4-trimethyl-	C9H2	
5.0922	1000144-04-3	7-Azabicyclo[4,2,0]octan-8-one	C6H1	
5.1030	106-65-0	Butanedioic acid, dimethyl ester	C12F	
5.3157	2801-84-5	Decane, 2,4-dimethyl-	C10H	
5.3618	7154-80-5	Heptane, 3,3,5-trimethyl-		1003
5.5285	122775-09-1	Cyclopropanecarboxylic acid, 2-methoxy-, methyl ester, trans-	1	115NO3S2
5.6200	299929-13-8	3-(4-Methylbenzoyl)-2-thioxo-4-thiazdlyl 4-methylbenzoate		1004
5.6388	617-52-7	Butanedioic acid, methylene-, dimethyl ester	C10	
5.6829	17302-01-1	3-Ethyl-3-methylheptane		8N2O
	613-94-5	Benzoic acid, hydrazide		120038
		Sulfurous acid, isobutyl pentyl ester		11004
5.7714	1000424-96-9	trans-2-Methyl-2-butenedioic acid dimethylester	-	11206
	1080-00-8	Methyl(methyl-4-deoxy.beta.l-threo-hex-4-enopyranosid)uronate	-	11005
5.8901	1000190-19-0	3-Acetoxy-3-hydroxypropionic acid, methyl ester		H12O4
5.9500	1604-11-1	Butanedioic acid, methyl-, dimethyl ester		H8OS
6.1698	13679-75-9	1-(2-Thienyl)-1-propanone		
6.2151	927-83-3	Diazene, bis(1,1-dimethylethyl)-	-	H18N2
6.2479	17865-32-6	Silane, cyclohexyldimethoxymethyl-		H20O2Si
6.4051	101-41-7	Benzeneacetic acid, methyl ester		H10O2
6.6228	18243-21-5	Formic acid, TMS derivative	C	H1002Si
6.6830	927-83-3	Diazene, bis(1,1-dimethylethyl)-	C	3H18N2
6.7617	20278-87-9	Heptane, 3,3,4-trimethyl-	C	10H22
6.8592	15764-16-6	Benzaldehyde, 2,4-dimethyl-	C	9H10O
and the same of the same of the same of	7154-80-5	Heptane, 3,3,5-trimethyl-	C	10H22
	7154-80-5	Heptane, 3,3,5-trimethyl-	C	10H22
	563-16-6	Hexane, 3,3-dimethyl-	C	28H18
		Undecane, 4,7-dimethyl-	(C13H28
	17301-32-5	2(1H)-Pyridinone, 4-hydroxy-1,6-dimethyl-		C7H9NO2
	6052-75-1			Accessed Supplemental Statement
7.6214 1	000453-46-5	Phenylglyoxylic acid, 2-butyl ester		C12H14O3
7.8138 4	110-44-5	Octane, 3,3-dimethyl-		C10H22
8.1765 2	5081-39-4	Benzoic acid, 3,5-dimethyl-, methyl ester		C10H12O2
8.4352 1	000115-35-8	Phenol, 2,4-bis(1-methylethyl)-, acetate		C14H20O2
and the second second	754-62-7	2-Propenoic acid, 3-phenyl-, methyl ester, (E)-		C10H10O2
8.9941 4		2,5-Furandicarboxylic acid, dimethyl ester		C8H8O5
0.3341 4.	and the second s	N-Ethylpyrrolidine-2,2-dicarboxilyc acid, dimethyl ester		C10H17NO

Page 1 of 2

Component RT	CAS#	Compound Name	Formula
9.0106	1000308-76-7	Diethylene glycol, O,O-di(pivaloyl)-	C14H26O5
9.0161	1000406-78-0	Pimelic acid, 3-(2-methoxyethyl)heptyl propyl ester	C20H38O5
9.0805	1732-09-8	Octanedioic acid, dimethyl ester	C10H18O4
9.1156	4271-99-2	1-Propene-1,2,3-tricarboxylic acid, trimethyl ester, (E)-	C9H12O6
9.2570	7154-80-5	Heptane, 3,3,5-trimethyl-	C10H22
9.3327	1587-20-8	Citric acid, trimethyl ester	C9H14O7
9.6484	1000309-19-5	Sulfurous acid, dodecyl 2-ethylhexyl ester	C20H42O3S
9.7504	22913-02-6	Nonaneperoxoic acid, 1,1-dimethylethyl ester	C13H26O3
9.8504	166273-38-7	Pentanoic acid, 5-hydroxy-, 2,4-di-t-butylphenyl esters	C19H30O3
10.0876	607-66-9	2(1H)-Quinolinone, 4-methyl-	C10H9NO
10.1318	4292-19-7	Dodecane, 1-iodo-	C12H25I
10.1706	1732-10-1	Nonanedioic acid, dimethyl ester	C11H20O4
10.2885	27037-61-2	8-Acetoxy-2-methylquinoline	C12H11NO2
10.3863	14526-12-6	Phosphor(isothiocyanatido)thioic difluoride	CF2NPS2
10.5456	5453-67-8	2,6-Pyridinedicarboxylic acid, dimethyl ester	C9H9NO4
11.6983	78607-80-4	2,2-Dimethyl-propyl 2,2-dimethyl-propane-thiosulfinate	C10H22OS2
11.9666	1292-19-7	Dodecane, 1-iodo-	C12H25I
and the second s	78607-80-4	2,2-Dimethyl-propyl 2,2-dimethyl-propane-thiosulfinate	C10H22OS2
12.4163 7	and the second s	Phosphoric acid, tris(2-ethylhexyl) ester	C24H51O4P
	267650-23-7	Tetradecanoic acid, 10,13-dimethyl-, methyl ester	C17H34O2
	8607-80-4	2,2-Dimethyl-propyl 2,2-dimethyl-propane-thiosulfinate	C10H22OS2
		Phthalic acid, cyclobutyl tridecyl ester	C25H38O4

Scanned with CamScanner

Qualitative Analysis Report

Data Filename

Pb lead_15 days.D

Sample Name

Pb lead_15 days

Sample Type

Instrument Name

GCMS-ALS

Position **User Name** 53 GCMSMS-HP\Preetam

Acq Method

Plant extractSCAN_.M

IRM Calibration Status

Not Applicable

Acquired Time DA Method

17/07/2023 11:57:58 default.m

Comment

Expected Barcode

Dual Inj Vol

Sample Amount

TuneName

atunes, eiex. tune.xml

TunePath D:\MassHunter\GCMS\1\7000 TuneDateStamp

2023-07-07T08:35:03+05:30

MSFirmwareVersion

DSP: 7000.3503, qqqServer: OperatorName G.7000.057-RUN

GCMSMS-HP\Preetam

RunCompletedFlag

True

Acquisition SW

Version

MassHunter GC/MS

Acquisition B.07.05.2479 23-Aug-2016 Copyright © 1989-

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Inc.

User Chromatograms

Fragmentor Voltage	Collision Energy	0	Ionizatio	n Mode	EI			_
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2111		•			,			1
1.75		1	-1					. 1
1.5		:						
1.25								
1 - 1022007914.60								
0.75								
0.5								
).25				٠.				
0						* :		
0 0 10 10 14 16	18 20 22 vs. Acquisition T	2 2 2 ime (r	4 26 nin)	28	30	32	34	36

Integration Peak List

Peak	Start	RT	End	Height	Area	Area %
1	3.072	3.095	3.432	50799872	353539598.2	2.57
2	3.592		3.901	1360024101	13754836569	100
3	5.047	5.073	5.138	213757602.4	287266240.4	2.09
4	5.276	5.306	5.337	108029109.8	149095843.1	1.08
5	6.81	6.85	6.923	632343469.4	1022007915	7.43

--- End Of Report ---

Qualitative Analysis Report

Data Filename

Pb lead_30 days.D

Sample Name

Position

Pb lead_30 days

Sample Type

Instrument Name Acq Method

GCMS-ALS Plant extractSCAN_.M **User Name Acquired Time** GCMSMS-HP\Preetam 17/07/2023 12:40:05

DA Method IRM Calibration Status Not Applicable

Comment

Expected Barcode

2 Dual Inj Vol **TunePath**

Sample Amount

TuneName

D:\MassHunter\GCMS\1\7000 TuneDateStamp

atunes.eiex.tune.xml

2023-07-07T08:35:03+05:30

default.m

MSFirmwareVersion

DSP: 7000.3503, qqqServer: OperatorName

GCMSMS-HP\Preetam

True RunCompletedFlag

Ģ.7000.057-RUN

Acquisition SW

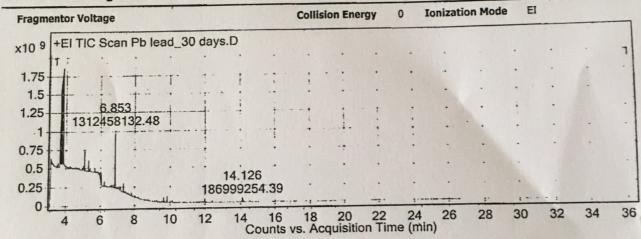
Version

MassHunter GC/MS

Acquisition B.07.05.2479 23-Aug-2016 Copyright © 1989-2016 Agilent Technologies,

Inc.

User Chromatograms



Peak	Start	RT	End	Height	Area	Area %
1	3.069	3.082	3.212	74252416	274621577.5	. 2.33
2	3.61			1292206848	11775596029	100
3	5.05		5.134	280328303.8	380503679.7	3.23
4	5.276		5.339	145865429.6	206149690.2	1.75
5	5.998			133576815.9	522885386.3	4.44
6	6.815	6.853	6.926	787953541.7	1312458132	11.15
7	7.322	7.347	7.401	118796780.1	165612431.5	1.41
8	9.604	9.647	9.691	65793301,06	120861757.7	1.03
9	9.808	9.845	9.877	81134722.61	120629748.	1.07
10	14.092	14.126		59277223.27	186999254.	4 1.5

--- End Of Report ---