

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU052120\
 Data File : VU038261.D
 Acq On : 21 May 2020 13:58
 Operator : JC/MD
 Sample : L2724-06DL 5X
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BE4T3DL

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.613	53	85	92	rBV	224959	559646	13.13%	2.032%
2	1.932	178	184	204	rVB2	169412	264058	6.20%	0.959%
3	2.401	320	330	350	rVB	756053	1168190	27.41%	4.242%
4	2.594	377	390	404	rBV	243906	399831	9.38%	1.452%
5	3.079	530	541	545	rBV3	35028	64182	1.51%	0.233%
6	3.112	547	551	565	rVB2	43801	74854	1.76%	0.272%
7	3.394	624	639	661	rVB2	95715	218715	5.13%	0.794%
8	3.735	730	745	765	rVB3	54803	125402	2.94%	0.455%
9	4.034	818	838	861	rBV	1245337	3244175	76.11%	11.780%
10	4.571	986	1005	1021	rBV	266021	646732	15.17%	2.348%
11	4.681	1021	1039	1085	rVB2	298884	976072	22.90%	3.544%
12	5.102	1154	1170	1202	rBV2	439852	1054723	24.74%	3.830%
13	5.423	1252	1270	1284	rBV3	37233	82552	1.94%	0.300%
14	5.526	1284	1302	1313	rBV3	23768	47779	1.12%	0.173%
15	5.761	1353	1375	1381	rBV2	517142	1288681	30.23%	4.679%
16	5.809	1382	1390	1406	rVB	907599	1808749	42.43%	6.568%
17	6.282	1522	1537	1569	rVB	456310	881730	20.69%	3.202%
18	6.722	1659	1674	1687	rBV	335075	648310	15.21%	2.354%
19	6.790	1687	1695	1706	rVB3	32232	59857	1.40%	0.217%
20	7.594	1931	1945	1960	rBV	161854	283567	6.65%	1.030%
21	7.742	1976	1991	2009	rVB3	41122	80073	1.88%	0.291%
22	7.925	2024	2048	2063	rBV	589821	1093330	25.65%	3.970%
23	7.992	2063	2069	2084	rVB2	38429	72765	1.71%	0.264%
24	8.205	2121	2135	2150	rBV	105305	177557	4.17%	0.645%
25	8.658	2262	2276	2313	rBV	1074941	2037630	47.80%	7.399%
26	8.815	2316	2325	2344	rVB2	21016	46825	1.10%	0.170%
27	9.468	2505	2528	2558	rVV2	1847235	4262427	100.00%	15.478%
28	9.713	2589	2604	2615	rBV2	30271	55565	1.30%	0.202%
29	9.780	2615	2625	2639	rBV2	27894	46360	1.09%	0.168%
30	10.121	2720	2731	2745	rVB5	48993	93111	2.18%	0.338%
31	10.500	2837	2849	2860	rBV	177954	282783	6.63%	1.027%
32	10.674	2897	2903	2913	rVB4	32379	52483	1.23%	0.191%
33	10.774	2914	2934	2960	rVB	299085	530617	12.45%	1.927%
34	10.912	2960	2977	2989	rBV5	189236	424082	9.95%	1.540%

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Integrator: RTE

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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 Title : TRACE VOA SOM01.0

35	11.047	2999	3019	3031	rBV2	55109	97923	2.30%	0.356%
36	11.307	3089	3100	3121	rVB	98692	183711	4.31%	0.667%
37	11.481	3145	3154	3164	rVB	47260	73235	1.72%	0.266%
38	11.655	3201	3208	3223	rVB8	48506	93403	2.19%	0.339%
39	11.825	3249	3261	3278	rVV2	716853	1448978	33.99%	5.262%
40	11.905	3278	3286	3306	rVB	107154	176895	4.15%	0.642%
41	12.108	3336	3349	3363	rVV2	257750	418382	9.82%	1.519%
42	12.208	3367	3380	3401	rVB	662191	1139589	26.74%	4.138%
43	12.587	3488	3498	3507	rBV	61843	94713	2.22%	0.344%
44	12.892	3583	3593	3610	rVB	133451	228709	5.37%	0.830%
45	12.995	3615	3625	3633	rVV	34228	55654	1.31%	0.202%
46	13.047	3633	3641	3652	rVB4	24369	44056	1.03%	0.160%
47	13.982	3923	3932	3939	rBV3	36215	61009	1.43%	0.222%
48	14.031	3939	3947	3959	rVB3	69168	117118	2.75%	0.425%
49	15.127	4272	4288	4303	rBV2	84700	152258	3.57%	0.553%

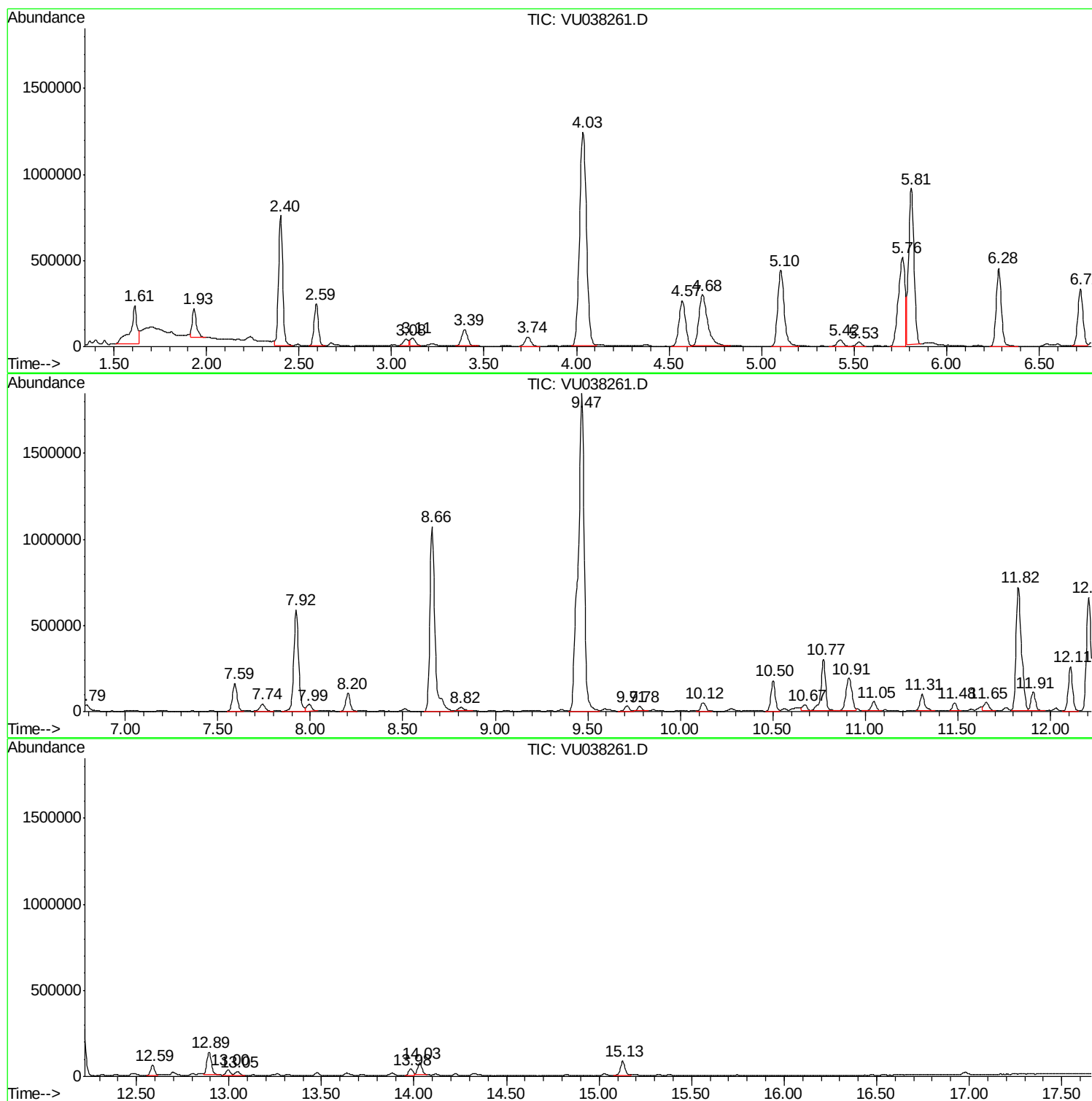
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Instrument :
MSVOA_U
ClientSampled :
BE4T3DL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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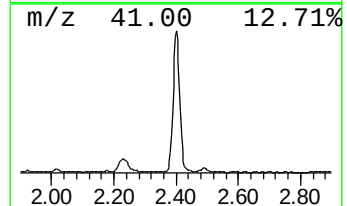
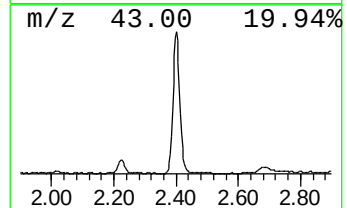
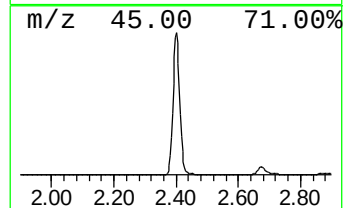
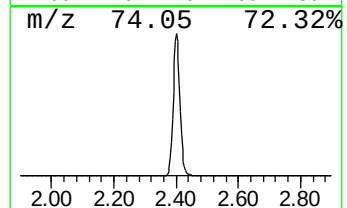
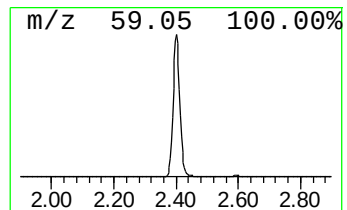
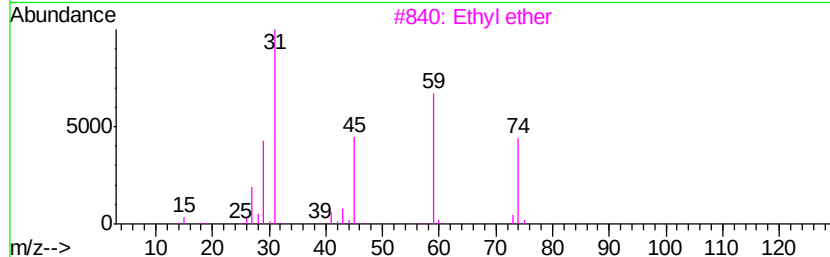
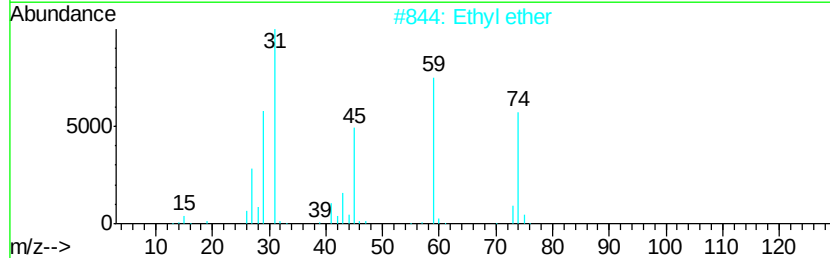
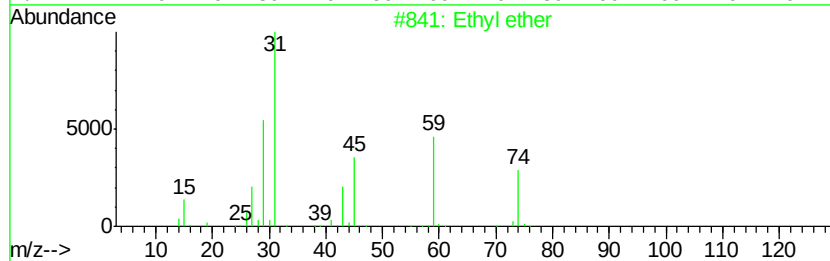
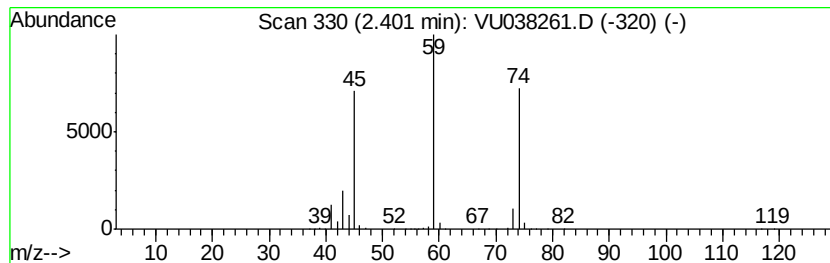
Instrument :
MSVOA_U
ClientSampleId :
BE4T3DL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethyl ether Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.40	6.62 ug/L	1168190	1,4-Difluorobenzene	6.28
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Ethyl ether	74 C4H10O	000060-29-7	90
2	Ethyl ether	74 C4H10O	000060-29-7	90
3	Ethyl ether	74 C4H10O	000060-29-7	90
4	Ethyl ether	74 C4H10O	000060-29-7	78
5	Ethane, 1,2-diethoxy-	118 C6H14O2	000629-14-1	56



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Instrument :
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ClientSampleId :
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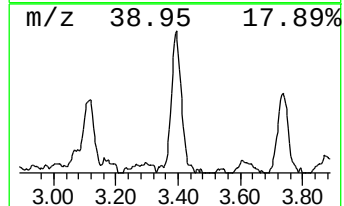
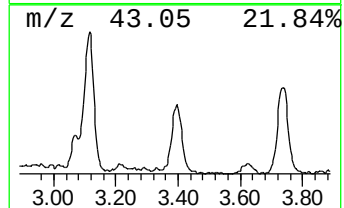
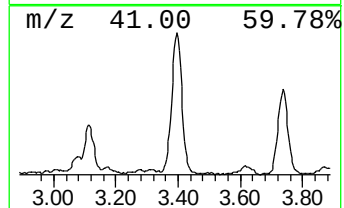
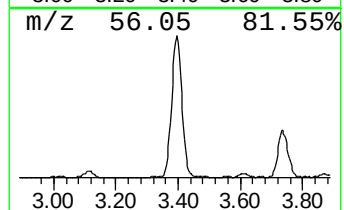
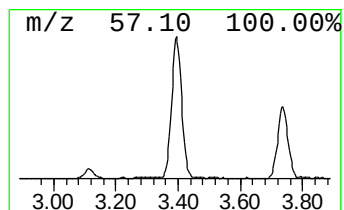
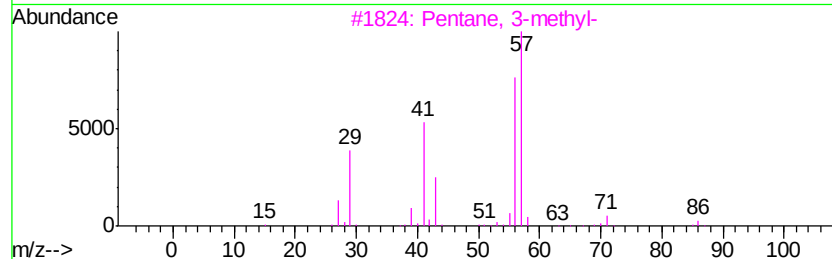
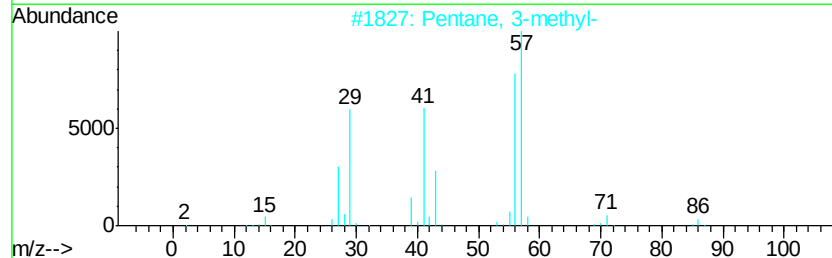
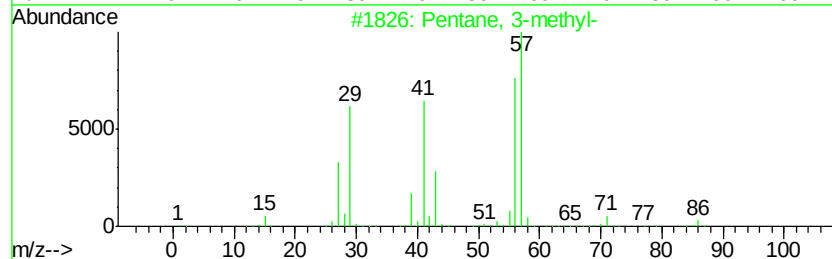
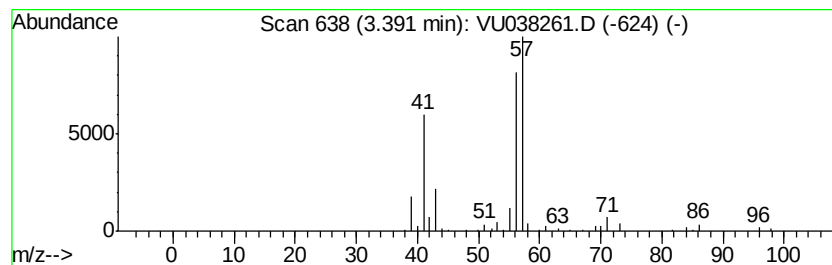
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.39	1.24 ug/L	218715	1,4-Difluorobenzene	6.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methyl-	86	C6H14	000096-14-0	80
2		Pentane, 3-methyl-	86	C6H14	000096-14-0	80
3		Pentane, 3-methyl-	86	C6H14	000096-14-0	80
4		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
5		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64



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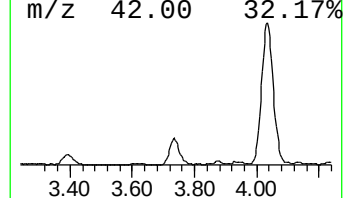
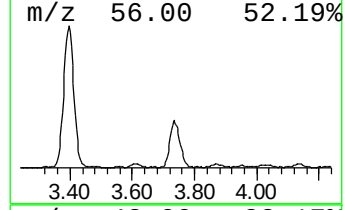
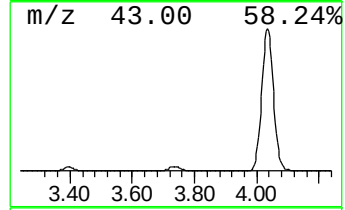
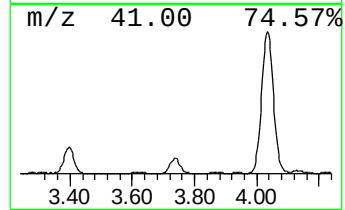
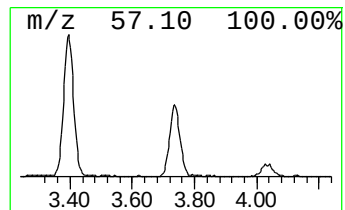
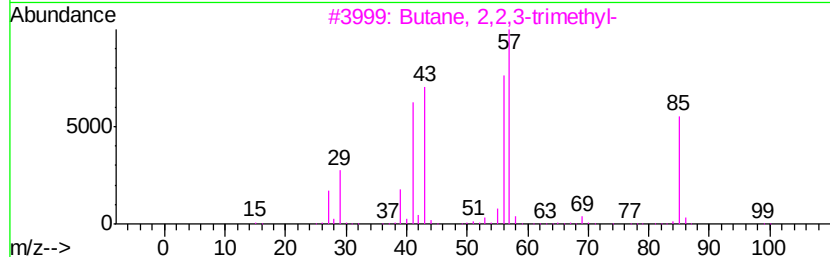
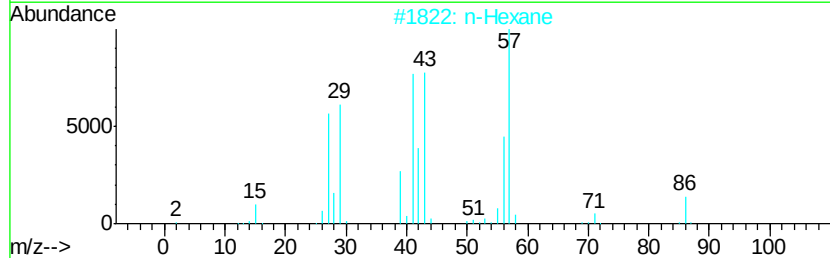
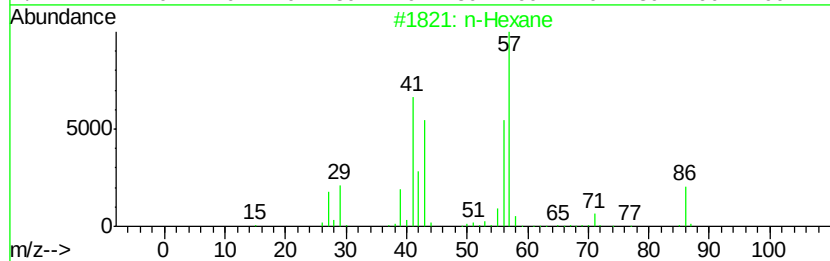
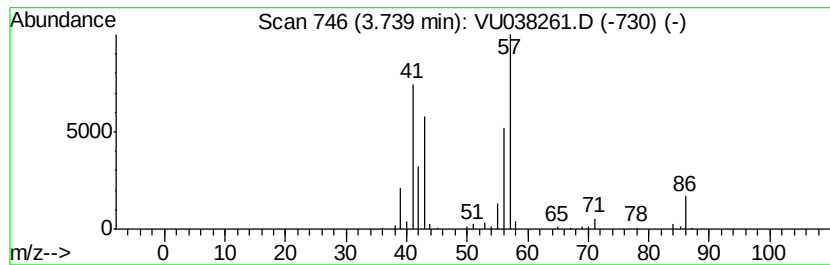
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
3.74	0.71 ug/L	125402	1,4-Difluorobenzene	6.28	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	91
2	n-Hexane	86	C6H14	000110-54-3	59
3	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	50
4	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	50
5	n-Hexane	86	C6H14	000110-54-3	50



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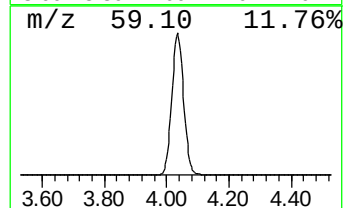
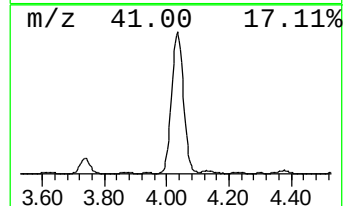
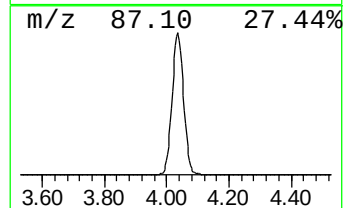
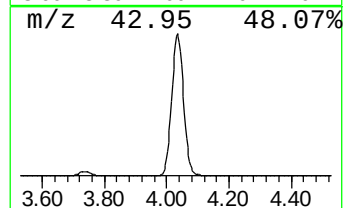
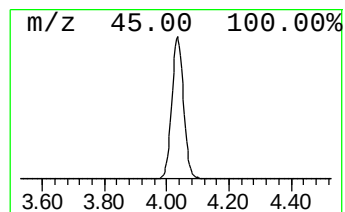
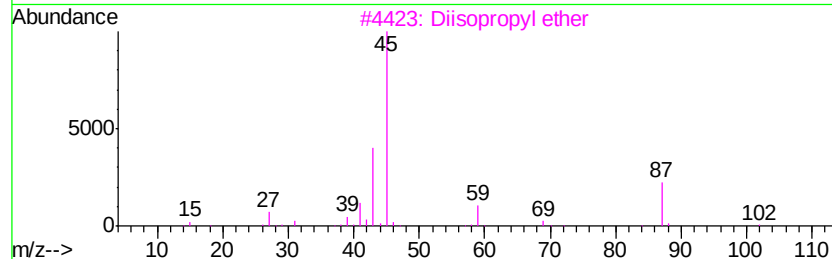
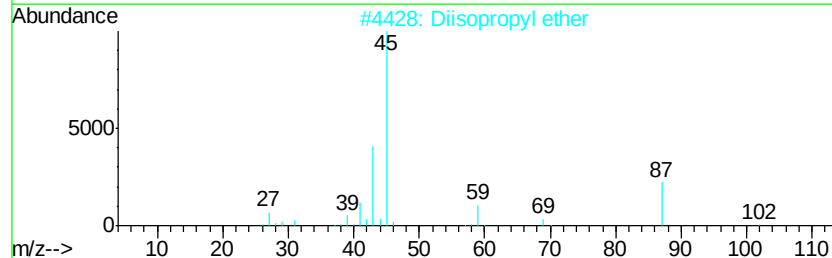
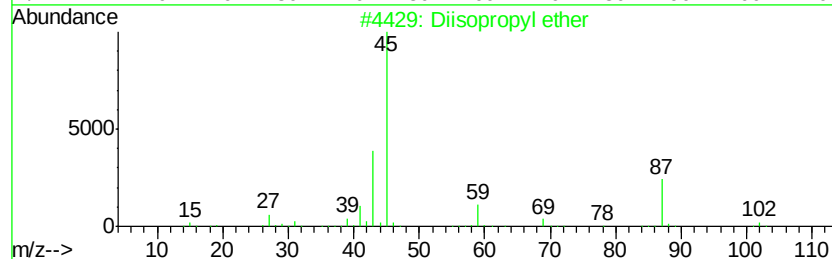
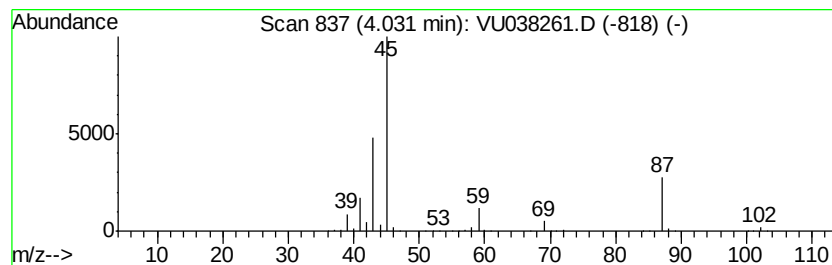
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Diisopropyl ether Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.03	18.40 ug/L	3244180	1,4-Difluorobenzene	6.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diisopropyl ether	102	C6H14O	000108-20-3	91
2		Diisopropyl ether	102	C6H14O	000108-20-3	83
3		Diisopropyl ether	102	C6H14O	000108-20-3	83
4		1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	74
5		Ether, 2-chloro-1-methylethyl is...	136	C6H13ClO	098277-76-0	64



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BE4T3DL

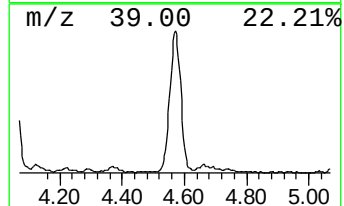
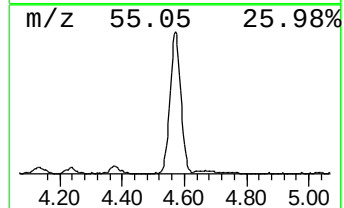
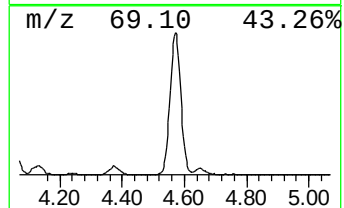
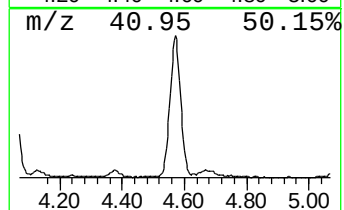
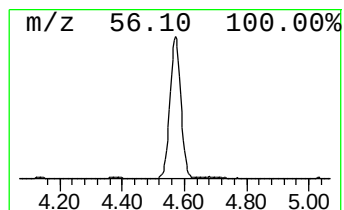
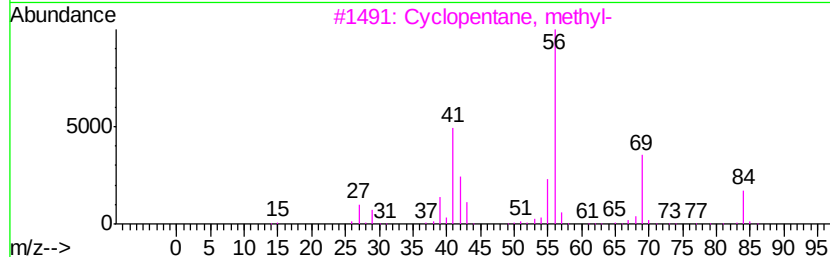
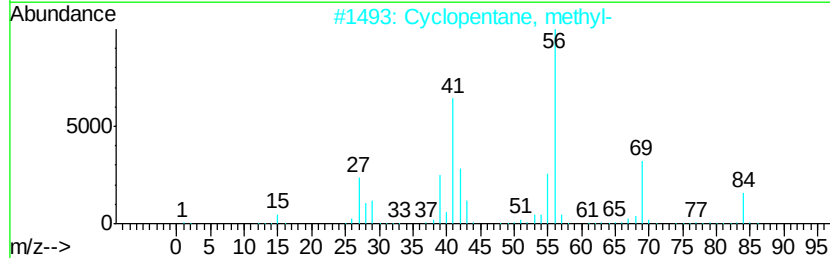
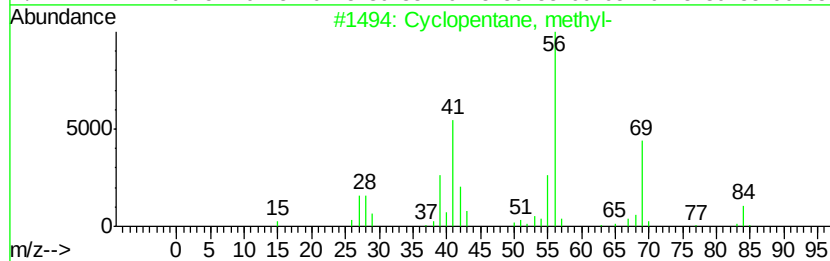
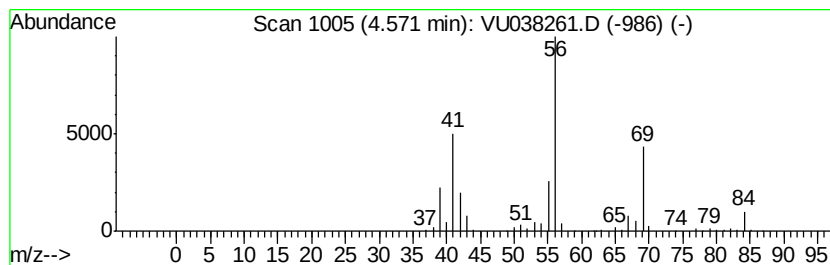
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Cyclic4.57 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.57	3.67 ug/L	646732	1,4-Difluorobenzene	6.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclopentane, methyl-	84	C6H12	000096-37-7	90
3		Cyclopentane, methyl-	84	C6H12	000096-37-7	78
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU052120\
Data File : VU038261.D
Acq On : 21 May 2020 13:58
Operator : JC/MD
Sample : L2724-06DL 5X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 5 Sample Multiplier: 1

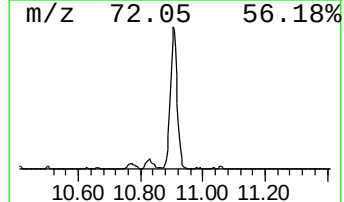
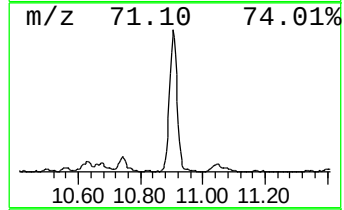
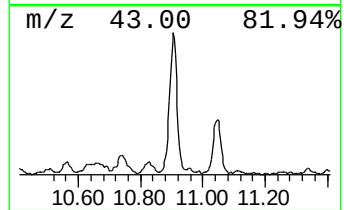
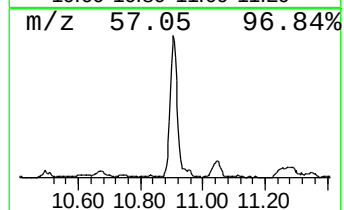
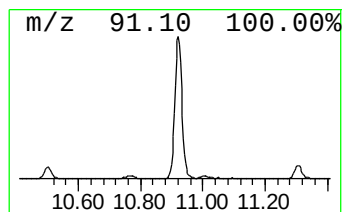
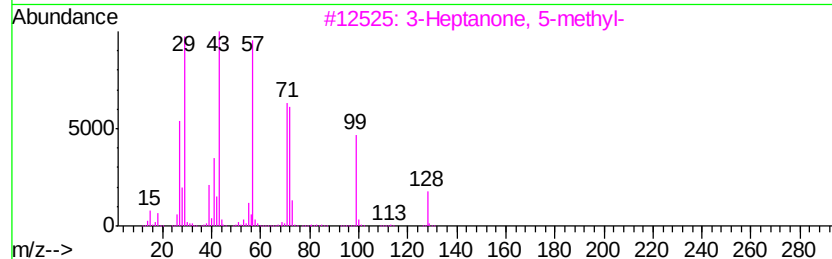
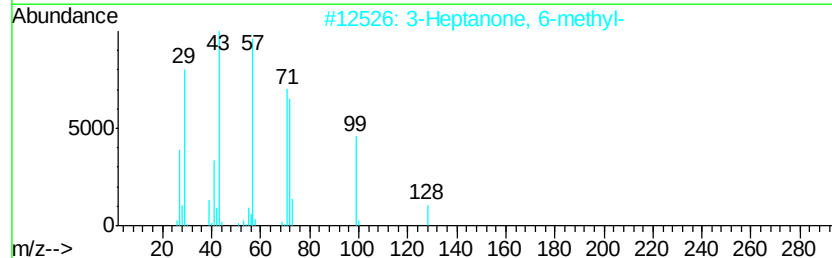
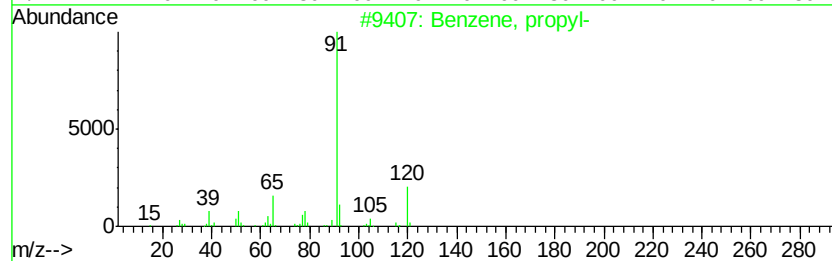
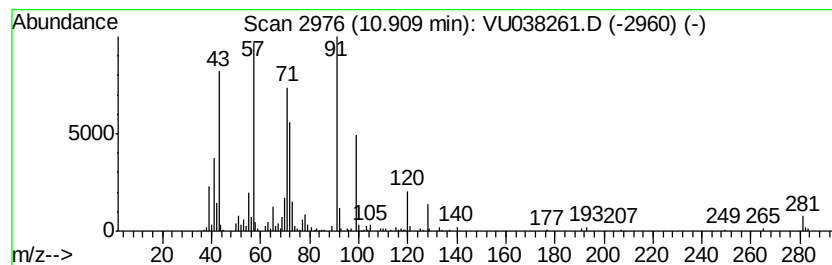
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, propyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.91	1.46 ug/L	424082	1,4-Dichlorobenzene-d4	11.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	64
2	3-Heptanone, 6-methyl-	128	C8H16O	000624-42-0	60
3	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	60
4	3-Octanone	128	C8H16O	000106-68-3	53
5	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU052120\
Data File : VU038261.D
Acq On : 21 May 2020 13:58
Operator : JC/MD
Sample : L2724-06DL 5X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 5 Sample Multiplier: 1

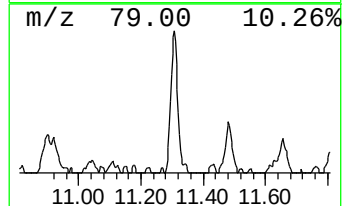
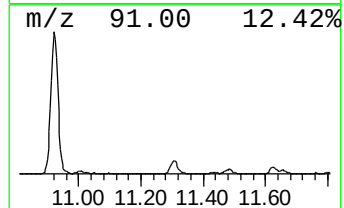
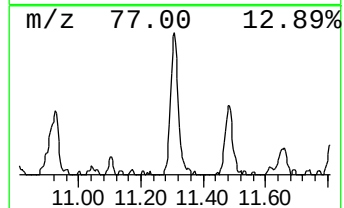
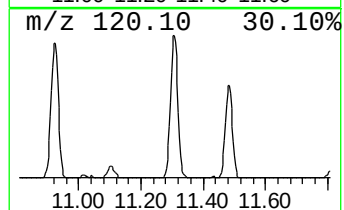
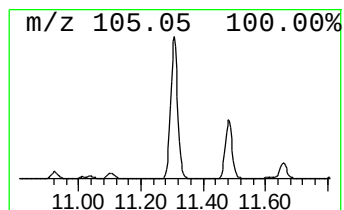
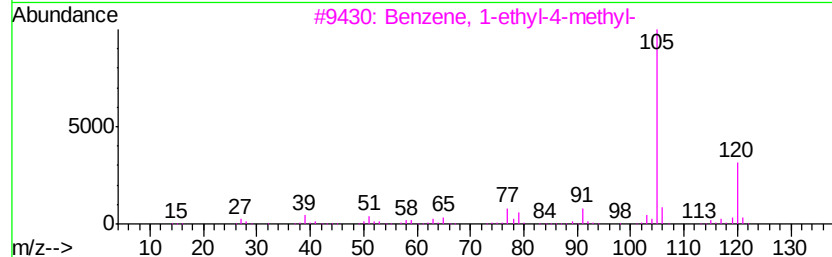
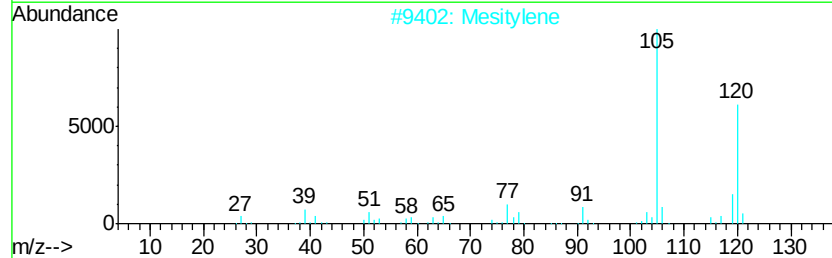
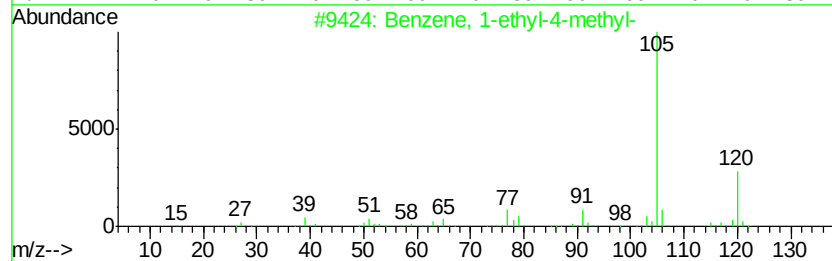
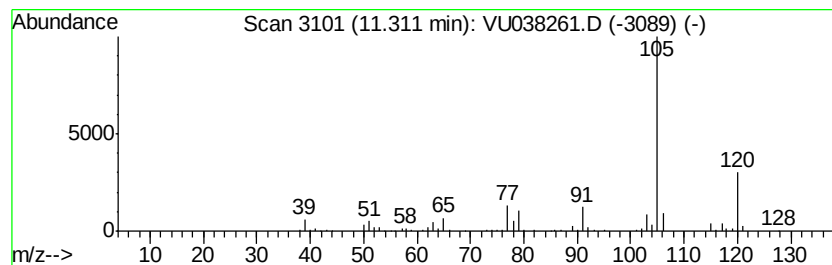
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1-ethyl-4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.31	0.63 ug/L	183711	1,4-Dichlorobenzene-d4	11.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 91
2		Mesitylene	120 C9H12	000108-67-8 91
3		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 91
4		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 91
5		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU052120\
Data File : VU038261.D
Acq On : 21 May 2020 13:58
Operator : JC/MD
Sample : L2724-06DL 5X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE4T3DL

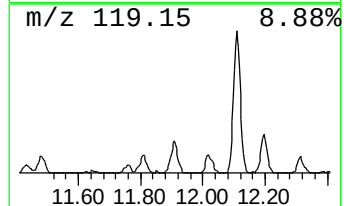
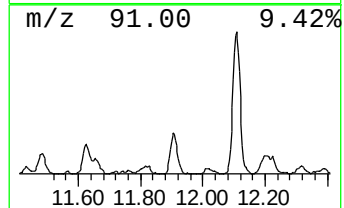
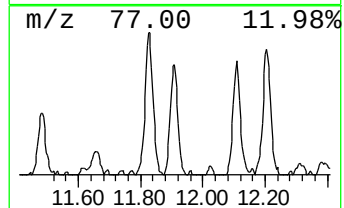
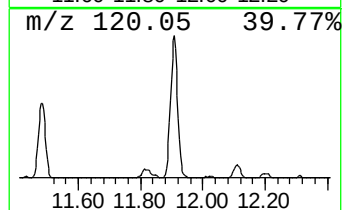
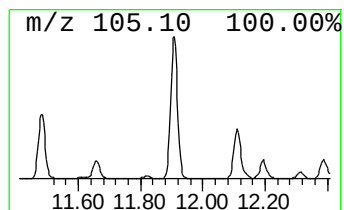
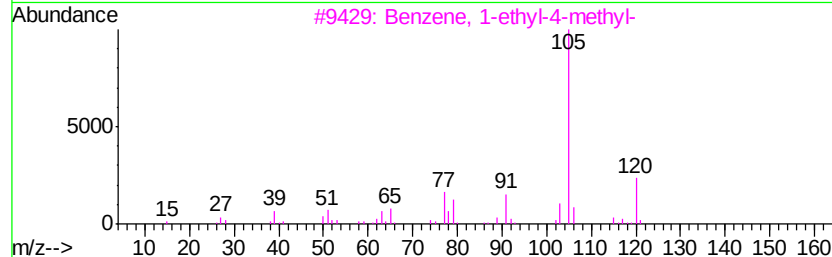
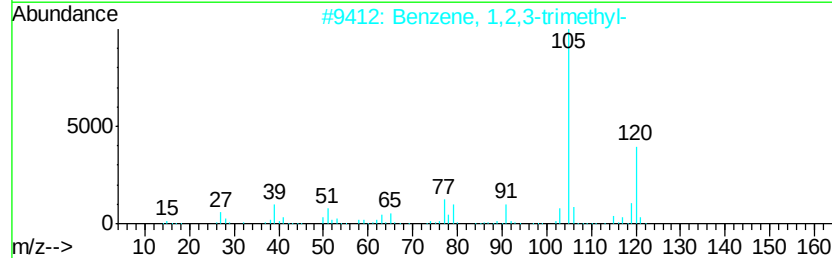
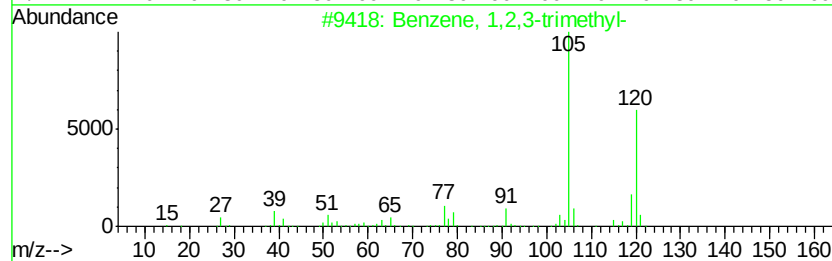
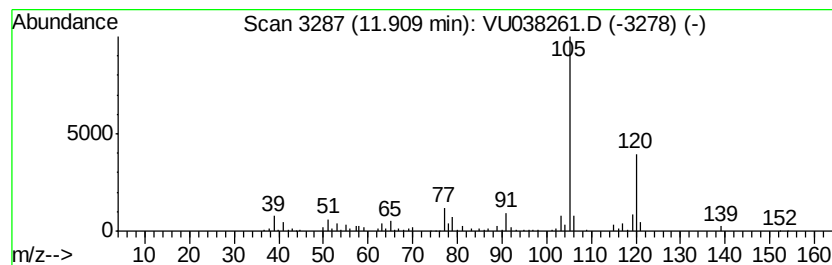
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.91	0.61 ug/L	176895	1,4-Dichlorobenzene-d4	11.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU052120\
Data File : VU038261.D
Acq On : 21 May 2020 13:58
Operator : JC/MD
Sample : L2724-06DL 5X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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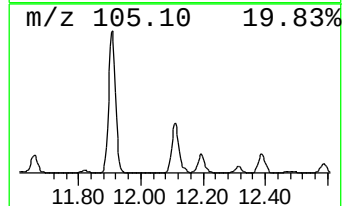
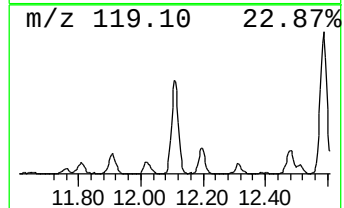
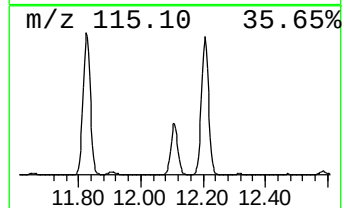
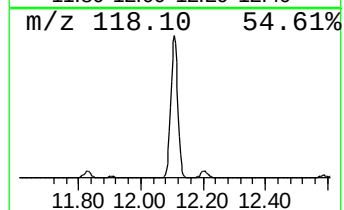
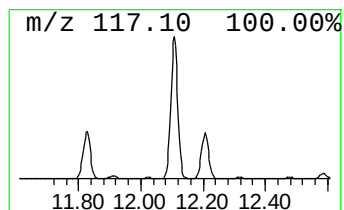
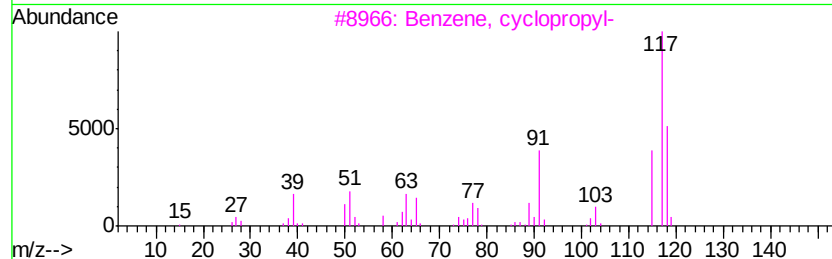
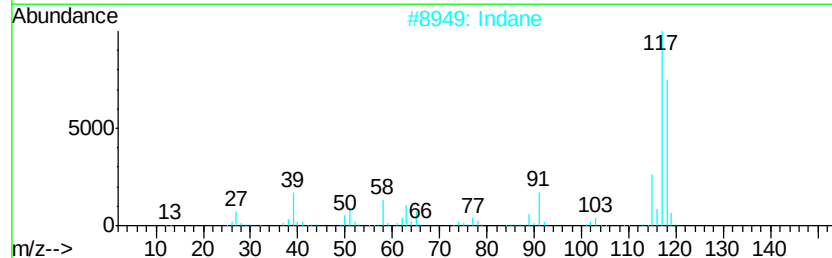
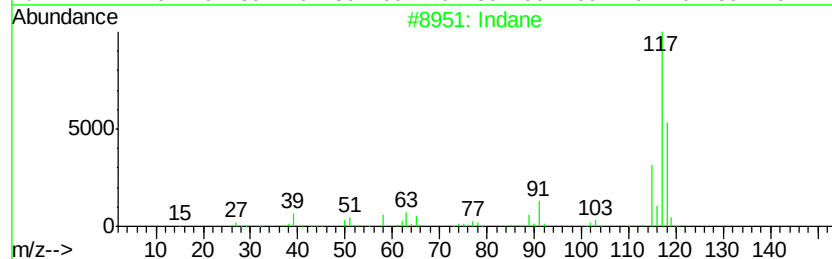
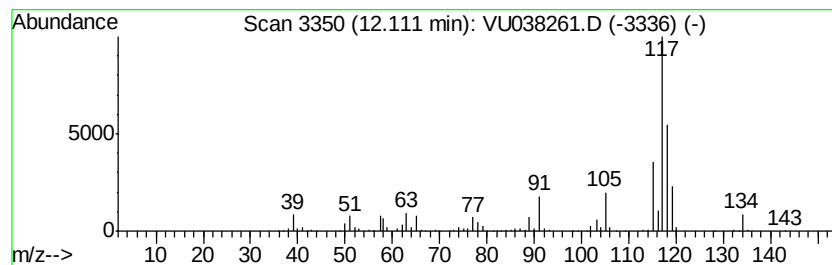
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.11	1.44 ug/L	418382	1,4-Dichlorobenzene-d4	11.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	93
2	Indane		118	C9H10	000496-11-7	76
3	Benzene, cyclopropyl-		118	C9H10	000873-49-4	70
4	Benzene, 2-propenyl-		118	C9H10	000300-57-2	68
5	Benzene, cyclopropyl-		118	C9H10	000873-49-4	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU052120\
Data File : VU038261.D
Acq On : 21 May 2020 13:58
Operator : JC/MD
Sample : L2724-06DL 5X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE4T3DL

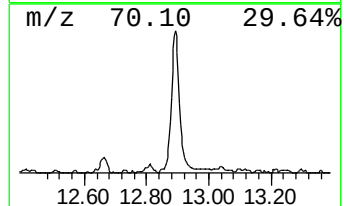
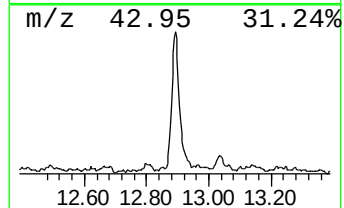
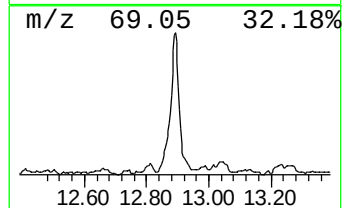
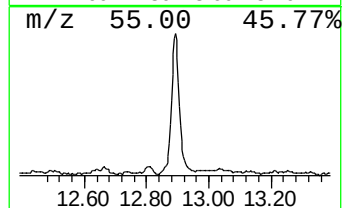
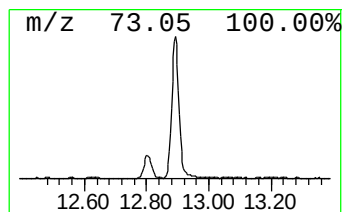
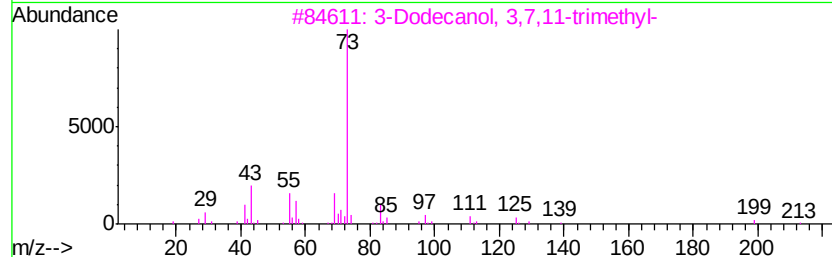
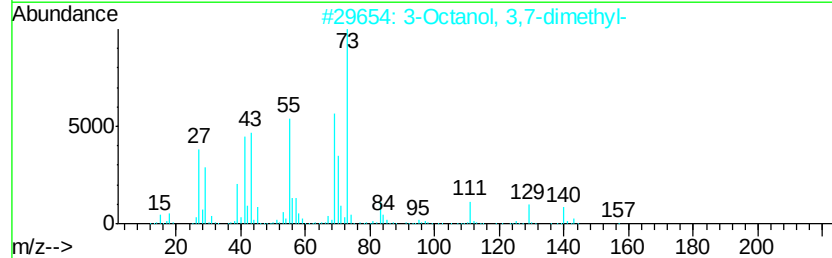
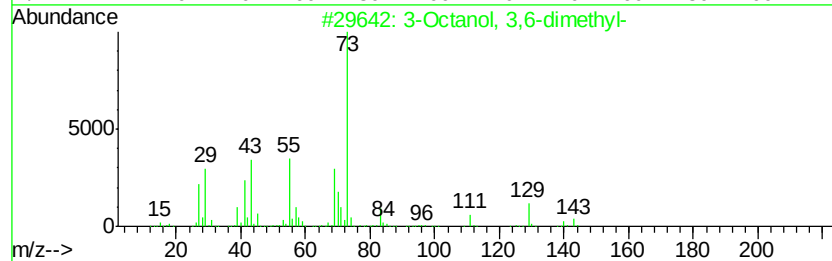
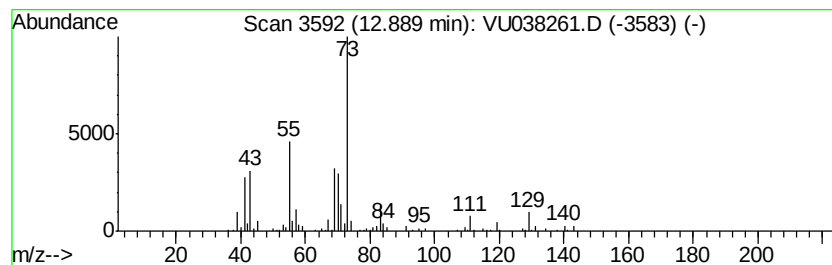
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 3-Octanol, 3,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.89	0.79 ug/L	228709	1,4-Dichlorobenzene-d4	11.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Octanol, 3,6-dimethyl-	158	C10H22O	000151-19-9	64
2		3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	43
3		3-Dodecanol, 3,7,11-trimethyl-	228	C15H32O	007278-65-1	42
4		3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	38
5		3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	28



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU052120\
Data File : VU038261.D
Acq On : 21 May 2020 13:58
Operator : JC/MD
Sample : L2724-06DL 5X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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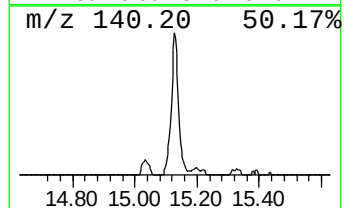
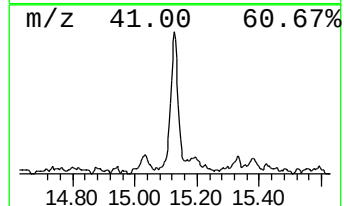
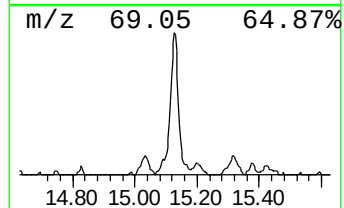
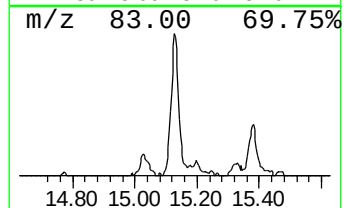
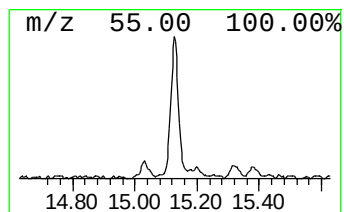
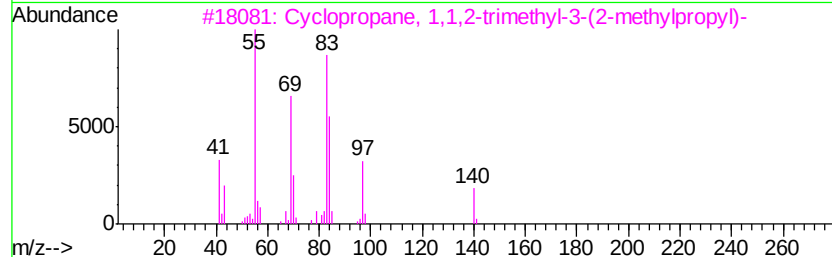
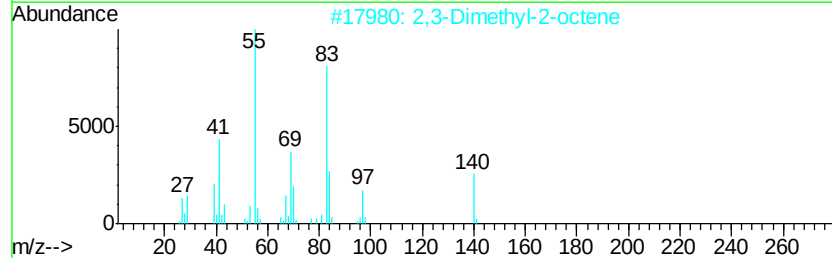
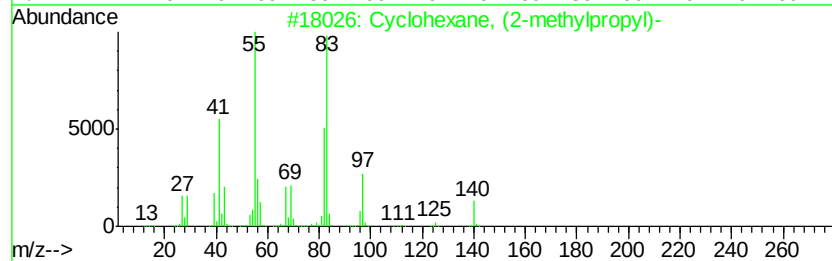
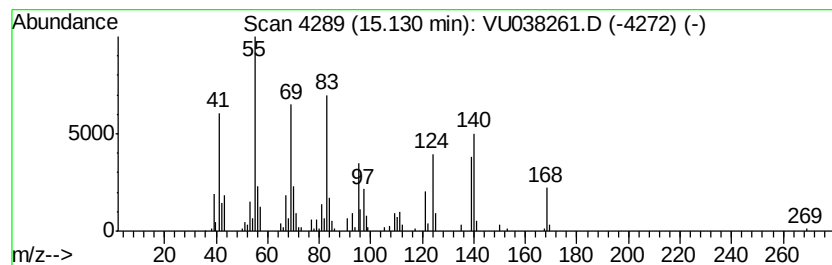
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 (DEL) Alkane: Cyclic15.13 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.13	0.53 ug/L	152258	1,4-Dichlorobenzene-d4	11.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	55
2		2,3-Dimethyl-2-octene	140	C10H20	019781-18-1	45
3		Cyclopropane, 1,1,2-trimethyl-3-...	140	C10H20	041977-43-9	38
4		Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	38
5		3-Dodecene, (Z)-	168	C12H24	007239-23-8	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU052120\
Data File : VU038261.D
Acq On : 21 May 2020 13:58
Operator : JC/MD
Sample : L2724-06DL 5X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE4T3DL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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(DEL) Alkane: Str...	3.39	1.2	ug/L	218715	1	6.28	881730	5.0
(DEL) Alkane: Str...	3.74	0.7	ug/L	125402	1	6.28	881730	5.0
Diisopropyl ether	4.03	18.4	ug/L	3244180	1	6.28	881730	5.0
(DEL) Alkane: Cyc...	4.57	3.7	ug/L	646732	1	6.28	881730	5.0
Benzene, propyl-	10.91	1.5	ug/L	424082	3	11.82	1448980	5.0
Benzene, 1-ethyl-...	11.31	0.6	ug/L	183711	3	11.82	1448980	5.0
Benzene, 1,2,3-tr...	11.91	0.6	ug/L	176895	3	11.82	1448980	5.0
Indane	12.11	1.4	ug/L	418382	3	11.82	1448980	5.0
3-Octanol, 3,6-di...	12.89	0.8	ug/L	228709	3	11.82	1448980	5.0
(DEL) Alkane: Cyc...	15.13	0.5	ug/L	152258	3	11.82	1448980	5.0