LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018753.D
Acq On : 02 Oct 2020 17:28
Operator : JC/SP
Sample : L4171-11
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BG3Q1

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.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	1.160	9	12	18	rBV2	72294	63167	7.75%	0.620%
2	1.240	22	25	27	rBV3	11400	9579	1.18%	0.094%
3	1.386	43	49	58	rBV2	125487	149541	18.35%	1.468%
4	1.483	61	65	69	rBV	34724	40865	5.01%	0.401%
5	1.697	95	100	106	rBV	83572	100180	12.29%	0.983%
6	1.935	136	139	144	rVB6	6594	9188	1.13%	0.090%
7	2.343	197	206	216	rBV	195577	312498	38.34%	3.067%
8	2.447	216	223	229	rBV2	17196	40150	4.93%	0.394%
9	2.550	235	240	245	rVB	10898	15869	1.95%	0.156%
10	2.831	280	286	288	rBV5	4712	9902	1.21%	0.097%
11	2.873	288	293	299	rVB3	16070	32628	4.00%	0.320%
12	3.148	328	338	348	rBV2	34671	85553	10.50%	0.840%
13	3.386	367	377	384	rVB8	6809	15838	1.94%	0.155%
14	3.501	386	396	406	rBV2	54328	124863	15.32%	1.225%
15	3.666	411	423	431	rBV4	7673	21535	2.64%	0.211%
16	4.300	515	527	532	rBV6	10718	35424	4.35%	0.348%
17	4.373	532	539	552	rVB3	30039	94434	11.58%	0.927%
18	4.544	557	567	580	rVV	98383	306399	37.59%	3.007%
19	4.660	580	586	593	rVB5	6631	20577	2.52%	0.202%
20	5.141	651	665	680	rBV2	124090	392387	48.14%	3.851%
21	6.044	800	813	826	rBV2	279936	815155	100.00%	8.000%
22	6.830	932	942	952	rBV	215550	508710	62.41%	4.992%
23	7.373	1021	1031	1041	rBV	165822	374879	45.99%	3.679%
24	7.635	1067	1074	1080	rBV6	9735	20953	2.57%	0.206%
25	8.372	1190	1195	1203	rVB	123493	192091	23.56%	1.885%
26	8.696	1242	1248	1254	rBV	449656	695135	85.28%	6.822%
27	8.763	1254	1259	1267	rVB	318400	501584	61.53%	4.922%
28	8.994	1288	1297	1303	rBV	83076	130797	16.05%	1.284%
29	9.348	1350	1355	1362	rVB3	7373	12780	1.57%	0.125%
30	9.427	1362	1368	1374	rBV	567065	780506	95.75%	7.660%
31	9.476	1374	1376	1383	rVB	13468	17930	2.20%	0.176%
32	9.561	1386	1390	1394	rBV2	22820	29816	3.66%	0.293%
33	10.098	1472	1478	1493	rBV	451141	627278	76.95%	6.156%
34	10.232	1493	1500	1505	rBV	24609	40289	4.94%	0.395%

LSC Area Percent Report

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Operator : JC/SP
Sample : L4171-11
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BG3Q1

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Title : TRACE VOA SOM01.0

35	10.342	1511	1518	1524	rBV3	42215	82593	10.13%	0.811%
36	10.396	1524	1527	1533	rVB3	24438	34487	4.23%	0.338%
37	10.500	1538	1544	1546	rBV6	6461	9726	1.19%	0.095%
38	10.628	1557	1565	1568	rBV4	26924	53839	6.60%	0.528%
39	10.683	1568	1574	1578	rVV	40384	67543	8.29%	0.663%
40	10.720	1578	1580	1585	rVV2	16946	19772	2.43%	0.194%
41	10.811	1589	1595	1600	rBV4	38932	74635	9.16%	0.732%
42	10.896	1605	1609	1614	rBV3	46811	76946	9.44%	0.755%
43	10.988	1620	1624	1625	rBV3	9576	11876	1.46%	0.117%
44	11.043	1630	1633	1636	rVV5	10150	16376	2.01%	0.161%
45	11.128	1640	1647	1649	rVV7	9167	20629	2.53%	0.202%
46	11.171	1651	1654	1658	rVV4	13527	18404	2.26%	0.181%
47	11.232	1658	1664	1675	rVV	189085	286639	35.16%	2.813%
48	11.341	1678	1682	1685	rVV	13573	19628	2.41%	0.193%
49	11.378	1685	1688	1690	rVV4	11643	15814	1.94%	0.155%
50	11.421	1691	1695	1700	rVV4	32744	70985	8.71%	0.697%
51	11.469	1700	1703	1706	rVV3	30105	47673	5.85%	0.468%
52	11.518	1708	1711	1715	rVV4	31812	41885	5.14%	0.411%
53	11.567	1715	1719	1722	rVV5	8328	13018	1.60%	0.128%
54	11.610	1722	1726	1729	rVV	89600	104998	12.88%	1.030%
55	11.652	1729	1733	1739	rVB5	30813	57958	7.11%	0.569%
56	11.793	1751	1756	1766	rVB	52984	89407	10.97%	0.877%
57	11.872	1766	1769	1771	rBV4	7034	9947	1.22%	0.098%
58	11.982	1781	1787	1790	rBV4	18742	33753	4.14%	0.331%
59	12.012	1790	1792	1795	rVB3	10149	9549	1.17%	0.094%
60	12.061	1795	1800	1806	rBV	518978	649197	79.64%	6.371%
61	12.128	1806	1811	1815	rBV	10801	14497	1.78%	0.142%
62	12.286	1834	1837	1839	rVV2	21906	29748	3.65%	0.292%
63	12.311	1839	1841	1844	rVV	21225	21776	2.67%	0.214%
64	12.360	1844	1849	1856	rVB	577040	769109	94.35%	7.548%
65	12.445	1858	1863	1868	rBV9	7485	15243	1.87%	0.150%
66	12.500	1868	1872	1876	rVB	18755	26453	3.25%	0.260%
67	12.573	1879	1884	1885	rBV	25677	30841	3.78%	0.303%
68	12.591	1885	1887	1892	rVB4	22428	30856	3.79%	0.303%
69	12.652	1892	1897	1903	rVB	46069	53409	6.55%	0.524%
70	12.756	1907	1914	1920	rBV9	12680	34382	4.22%	0.337%
71	12.890	1932	1936	1941	rVB3	48322	63768	7.82%	0.626%

LSC Area Percent Report

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ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BG3Q1

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Title : TRACE VOA SOM01.0

72	12.963	1941	1948	1951	rBV	50367	70181	8.61%	0.689%
73	13.000	1951	1954	1961	rVB	30870	43283	5.31%	0.425%
74	13.061	1961	1964	1966	rBV3	7021	9006	1.10%	0.088%
75	13.140	1974	1977	1981	rVB6	7317	10208	1.25%	0.100%
76	13.201	1981	1987	1993	rBV9	8287	17060	2.09%	0.167%
77	13.292	1997	2002	2004	rBV3	9592	15093	1.85%	0.148%
78	13.329	2004	2008	2014	rVV3	27847	48549	5.96%	0.476%
79	13.408	2018	2021	2023	rVB	8898	8714	1.07%	0.086%
80	13.451	2023	2028	2034	rBV10	10349	18075	2.22%	0.177%
81	13.621	2051	2056	2064	rVB5	26812	56558	6.94%	0.555%
82	13.725	2068	2073	2076	rBV4	14691	22313	2.74%	0.219%
83	13.817	2082	2088	2095	rVV2	43459	64582	7.92%	0.634%
84	13.890	2098	2100	2107	rVV8	9337	13447	1.65%	0.132%
85	13.969	2109	2113	2116	rVV4	7180	10299	1.26%	0.101%
86	14.006	2116	2119	2123	rVB6	7304	9091	1.12%	0.089%
87	14.109	2134	2136	2142	rVB7	7638	8709	1.07%	0.085%
88	14.256	2156	2160	2165	rVB7	6192	9442	1.16%	0.093%
89	14.359	2173	2177	2182	rBV6	6341	9346	1.15%	0.092%
90	14.481	2192	2197	2202	rBV9	5827	9789	1.20%	0.096%
91	14.676	2223	2229	2234	rBV	22561	33682	4.13%	0.331%
92	14.822	2248	2253	2256	rBV	11404	15773	1.93%	0.155%
93	15.530	2364	2369	2372	rBV7	8844	13783	1.69%	0.135%
94	15.664	2385	2391	2395	rBV6	10754	17101	2.10%	0.168%

Sum of corrected areas: 10189923

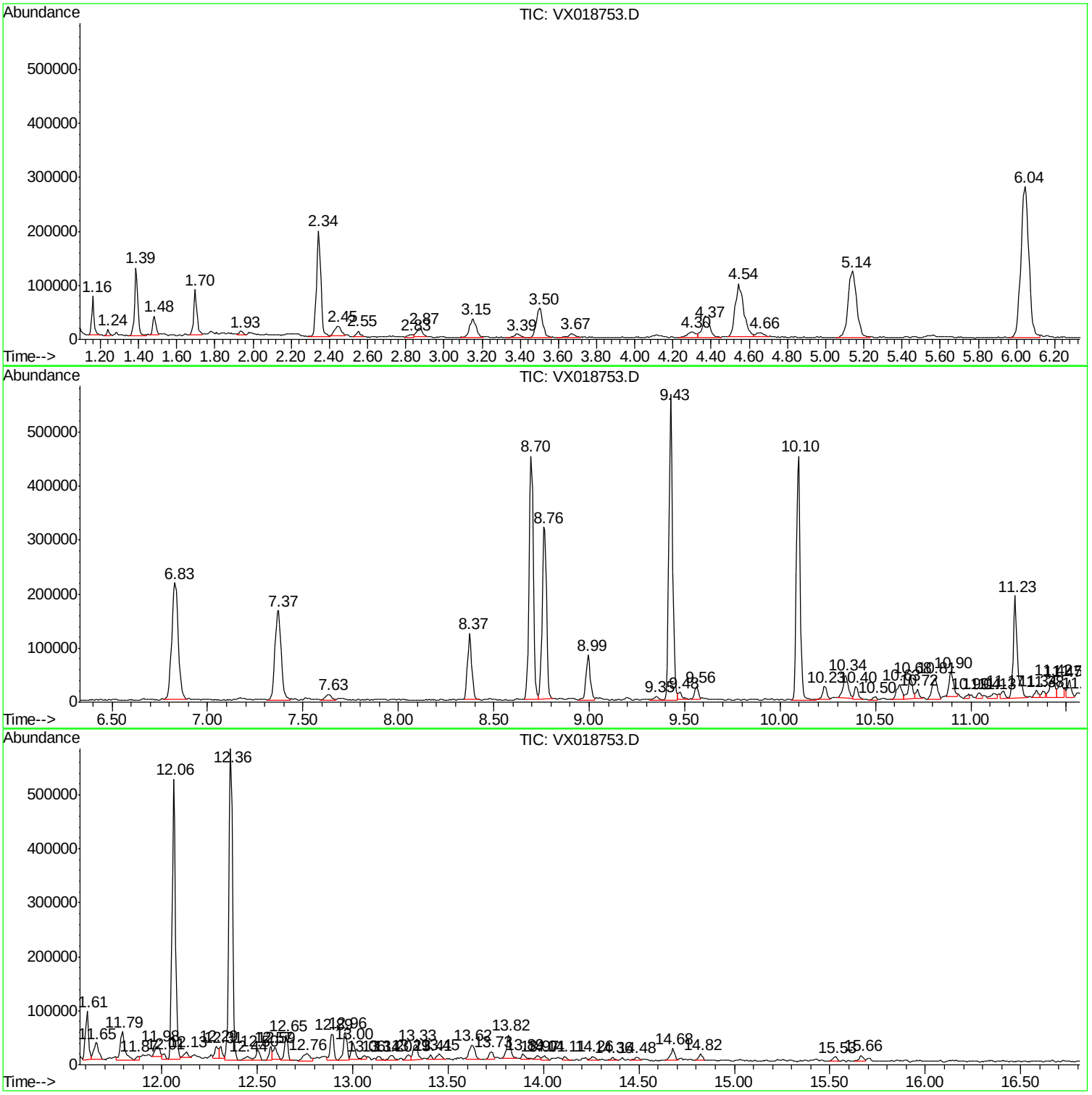
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018753.D
Acq On : 02 Oct 2020 17:28
Operator : JC/SP
Sample : L4171-11
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BG3Q1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018753.D
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Operator : JC/SP
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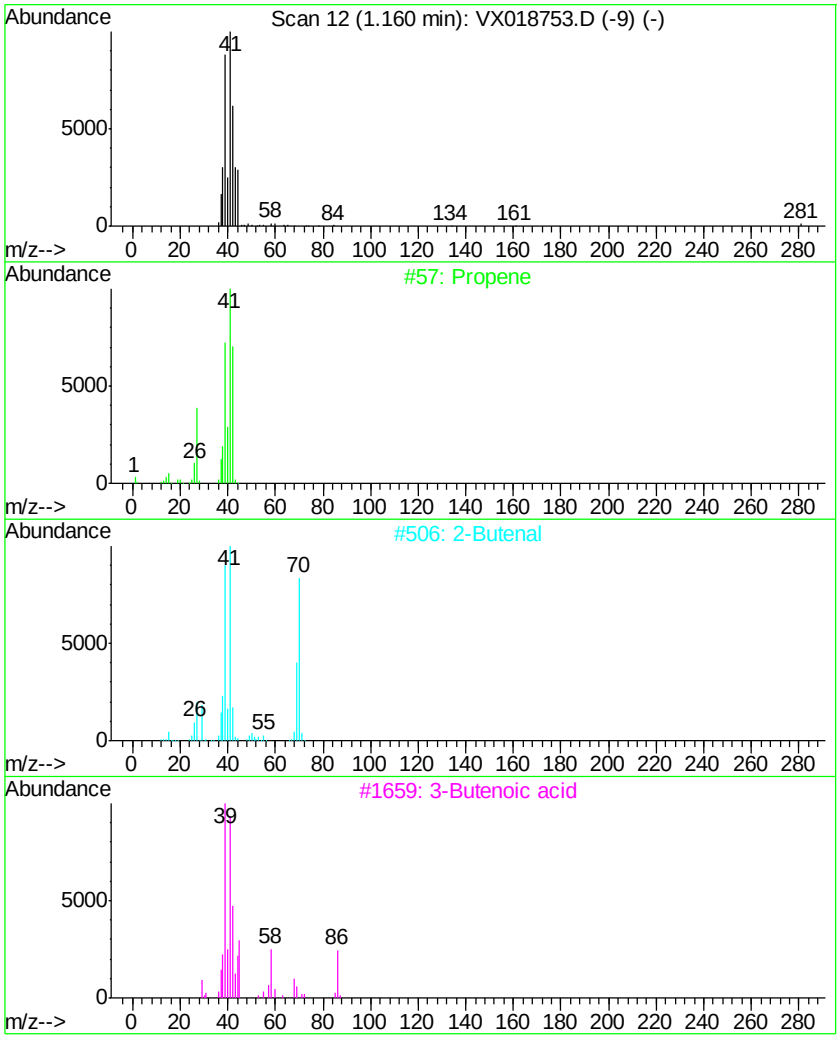
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

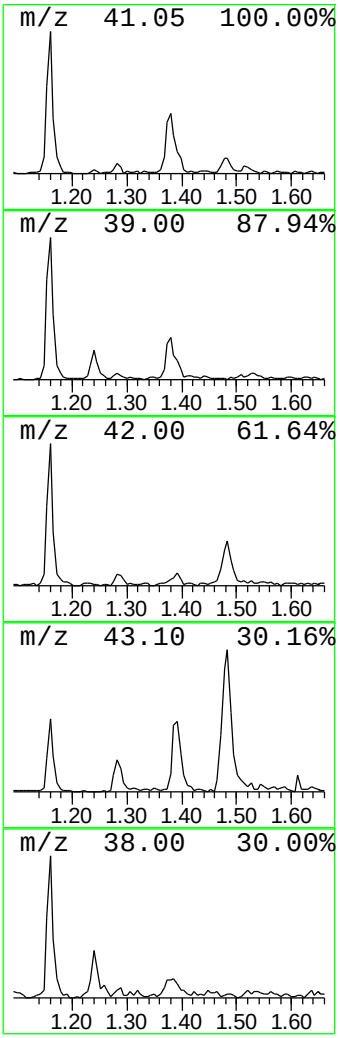
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.16	0.62 ug/L	63167	1,4-Difluorobenzene	6.83

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propene	42	C3H6	000115-07-1	72
2	2-Butenal	70	C4H6O	004170-30-3	64
3	3-Butenoic acid	86	C4H6O2	000625-38-7	64
4	2-Butenal, (E)-	70	C4H6O	000123-73-9	64
5	Cyclopropanecarboxylic acid	86	C4H6O2	001759-53-1	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018753.D
Acq On : 02 Oct 2020 17:28
Operator : JC/SP
Sample : L4171-11
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BG3Q1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

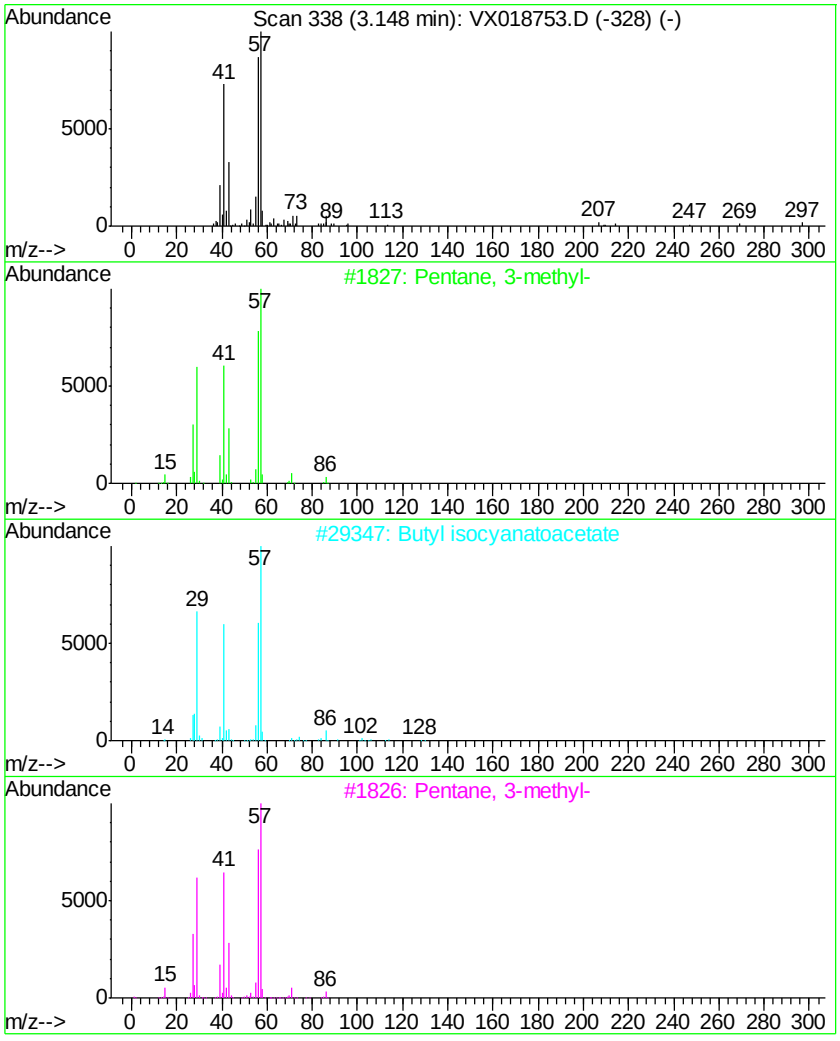
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

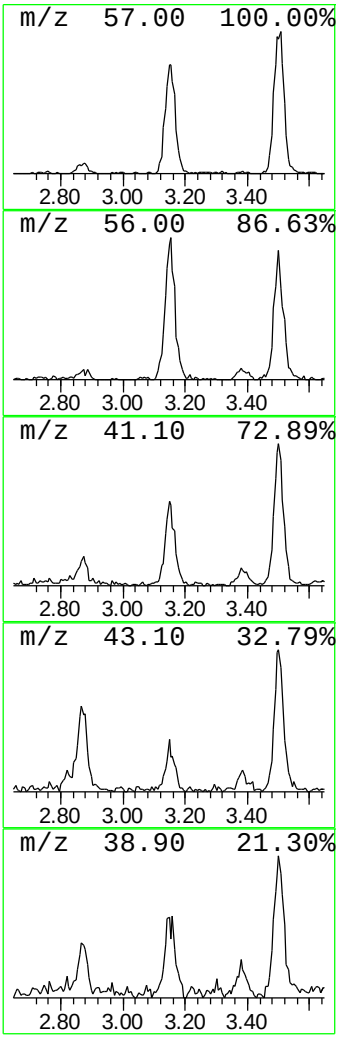
****

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	0.84 ug/L	85553	1,4-Difluorobenzene	6.83

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	72
2	Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	64
3	Pentane, 3-methyl-	86	C6H14	000096-14-0	64
4	Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	64
5	Pentane, 3-methyl-	86	C6H14	000096-14-0	50





Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018753.D
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Sample : L4171-11
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BG3Q1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

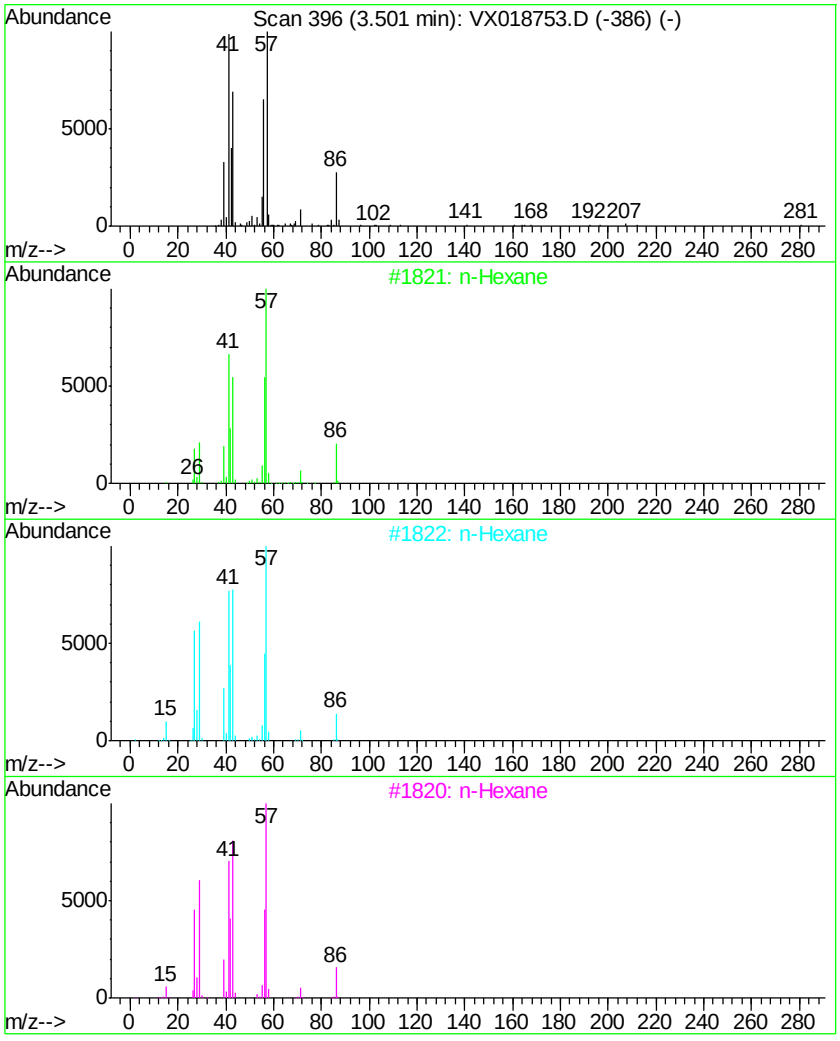
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

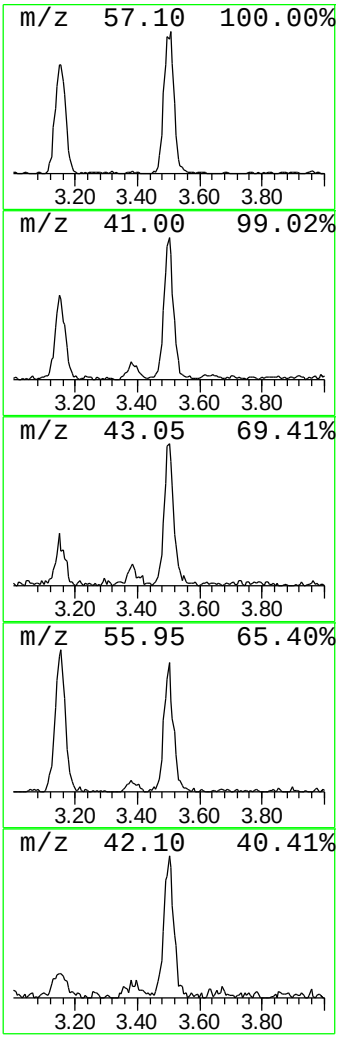
****

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.50	1.23 ug/L	124863	1,4-Difluorobenzene	6.83

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	76
2	n-Hexane	86	C6H14	000110-54-3	59
3	n-Hexane	86	C6H14	000110-54-3	59
4	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	40
5	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	40





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100220\
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Acq On : 02 Oct 2020 17:28
Operator : JC/SP
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Misc : 25.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BG3Q1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

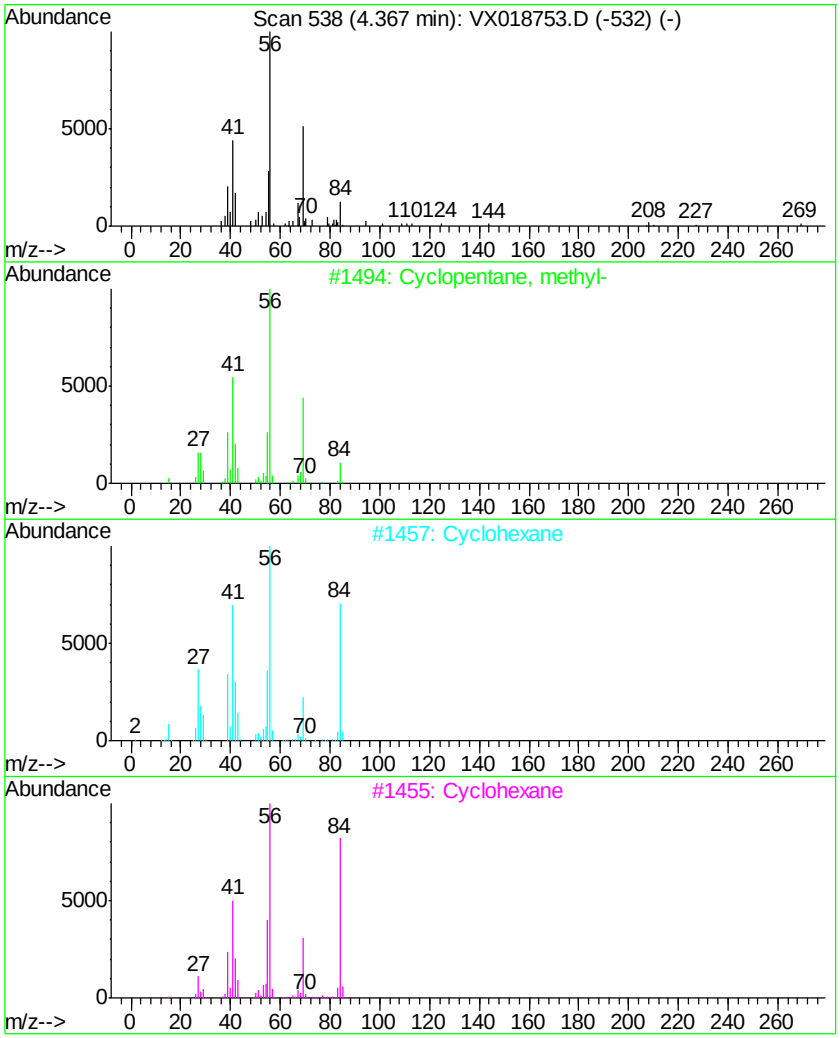
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

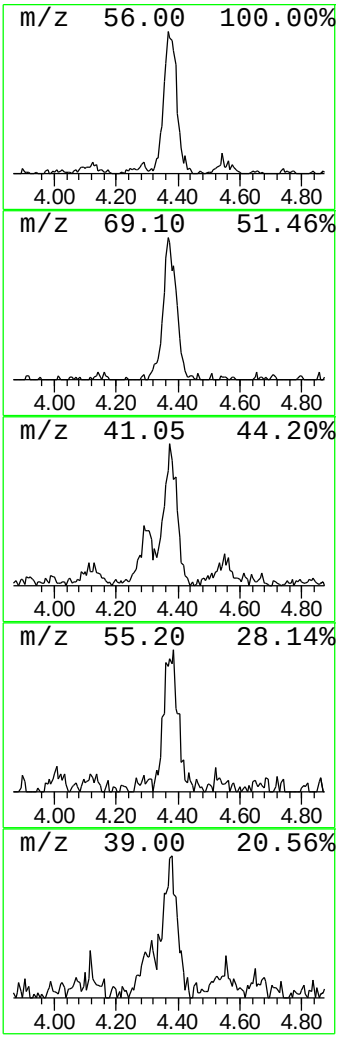
****

Peak Number 4 (DEL) Alkane: Cyclic4.37 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	0.93 ug/L	94434	1,4-Difluorobenzene	6.83

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	87
2	Cyclohexane	84	C6H12	000110-82-7	80
3	Cyclohexane	84	C6H12	000110-82-7	72
4	Cyclohexane	84	C6H12	000110-82-7	72
5	Hexyl trifluoroacetate	198	C8H13F3O2	000400-61-3	72





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Sample : L4171-11
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BG3Q1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

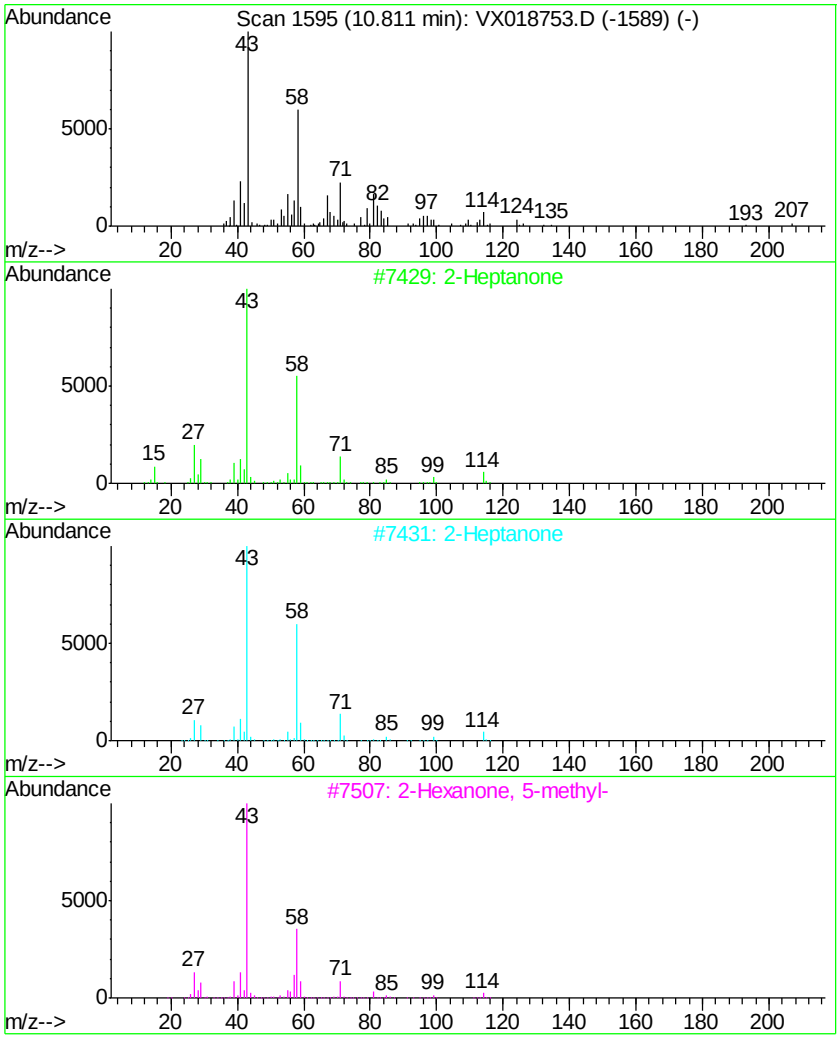
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

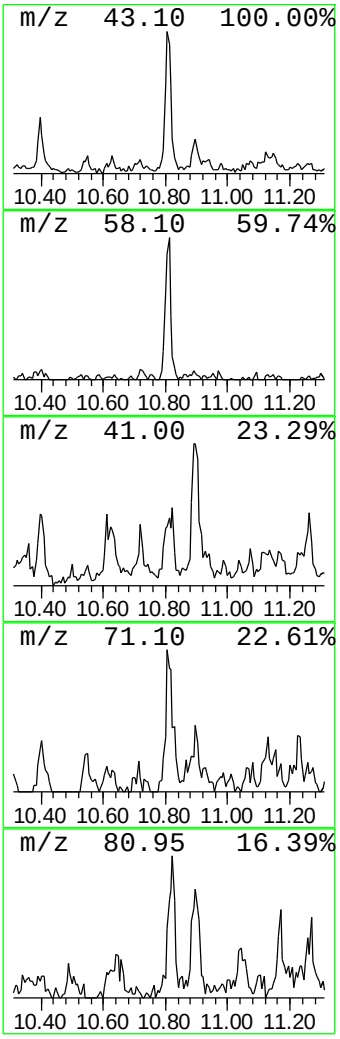
****

Peak Number 6 2-Heptanone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.81	0.59 ug/L	74635	Chlorobenzene-d5	10.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone	114	C7H14O	000110-43-0	58
2	2-Heptanone	114	C7H14O	000110-43-0	52
3	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	50
4	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	50
5	2-Heptanone	114	C7H14O	000110-43-0	50





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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

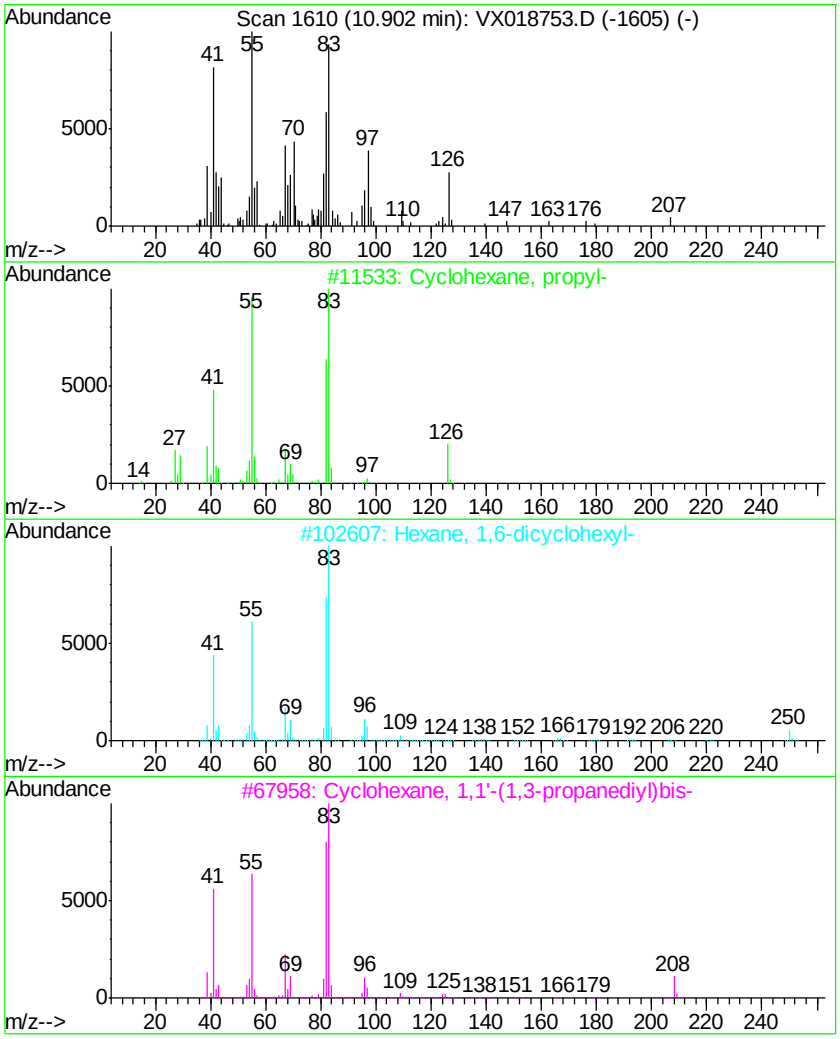
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

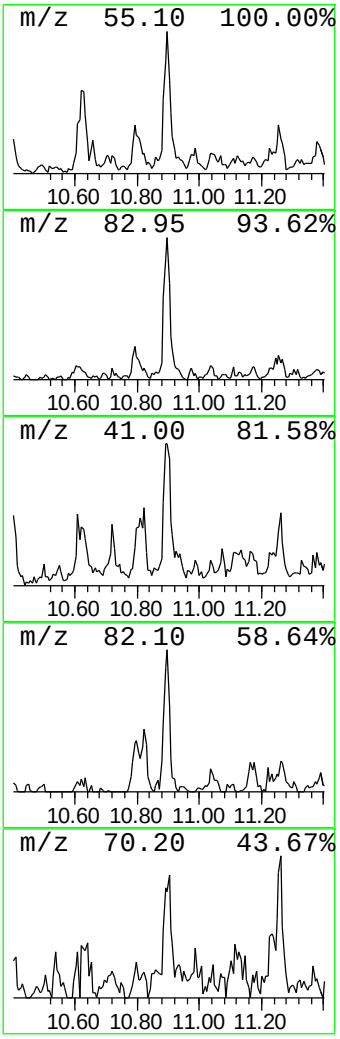
****

Peak Number 7 (DEL) Alkane: Cyclic10.90 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	0.61 ug/L	76946	Chlorobenzene-d5	10.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, propyl-	126	C9H18	001678-92-8	87
2	Hexane, 1,6-dicyclohexyl-	250	C18H34	001610-23-7	59
3	Cyclohexane, 1,1'-(1,3-propanedi...	208	C15H28	003178-24-3	59
4	Cyclohexane, 1,1'-(1,3-propanedi...	208	C15H28	003178-24-3	59
5	2,3-Dimethyl-2-heptene	126	C9H18	003074-64-4	52





Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018753.D
Acq On : 02 Oct 2020 17:28
Operator : JC/SP
Sample : L4171-11
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BG3Q1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

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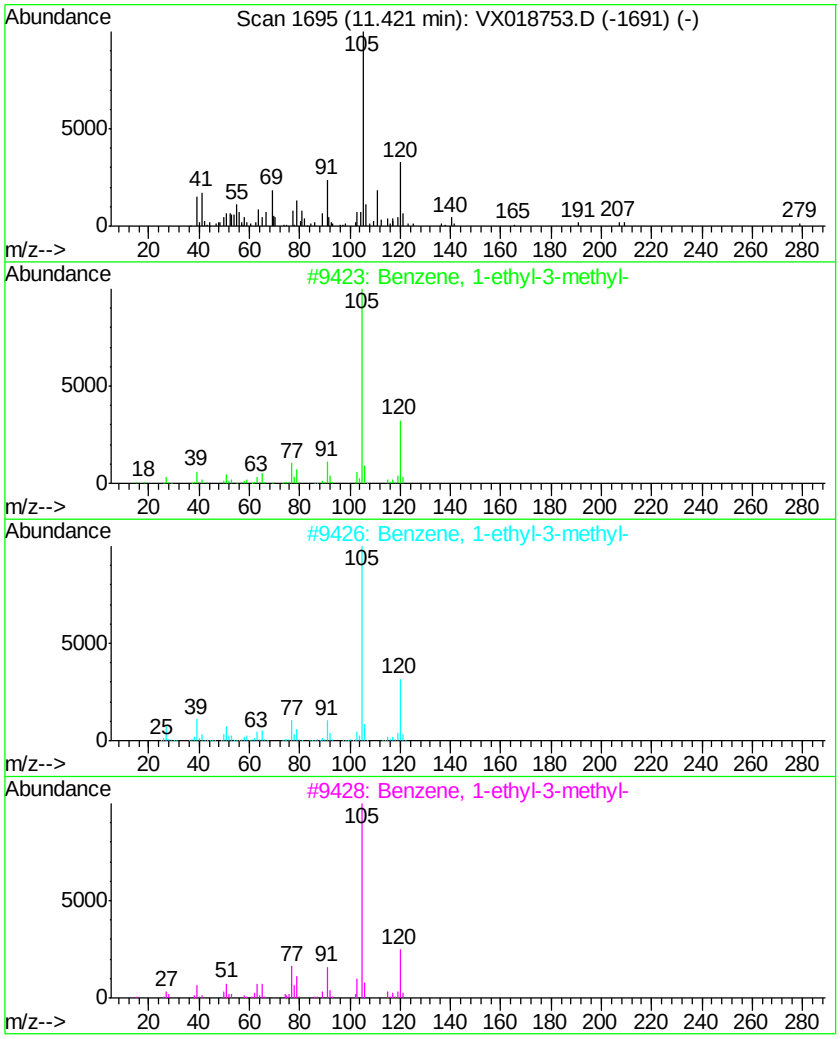
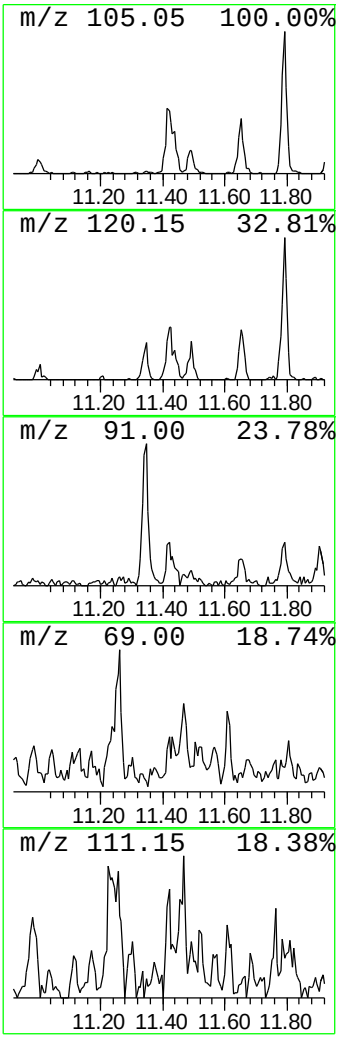
Peak Number 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	0.55 ug/L	70985	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	52
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	52
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	52
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	52
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	52





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100220\
Data File : VX018753.D
Acq On : 02 Oct 2020 17:28
Operator : JC/SP
Sample : L4171-11
Misc : 25.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BG3Q1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

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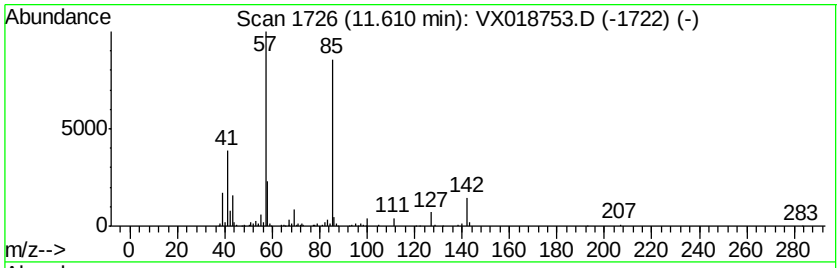
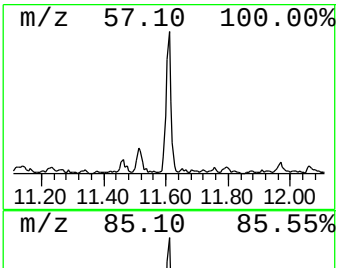
Peak Number 9 4-Heptanone, 2,6-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.61	0.81 ug/L	104998	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	91
2	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	72
3	5-Nonanone	142	C9H18O	000502-56-7	64
4	t-Butyl isobutyl ketone	142	C9H18O	014705-50-1	64
5	5-Nonanone	142	C9H18O	000502-56-7	59





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100220\
Data File : VX018753.D
Acq On : 02 Oct 2020 17:28
Operator : JC/SP
Sample : L4171-11
Misc : 25.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BG3Q1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

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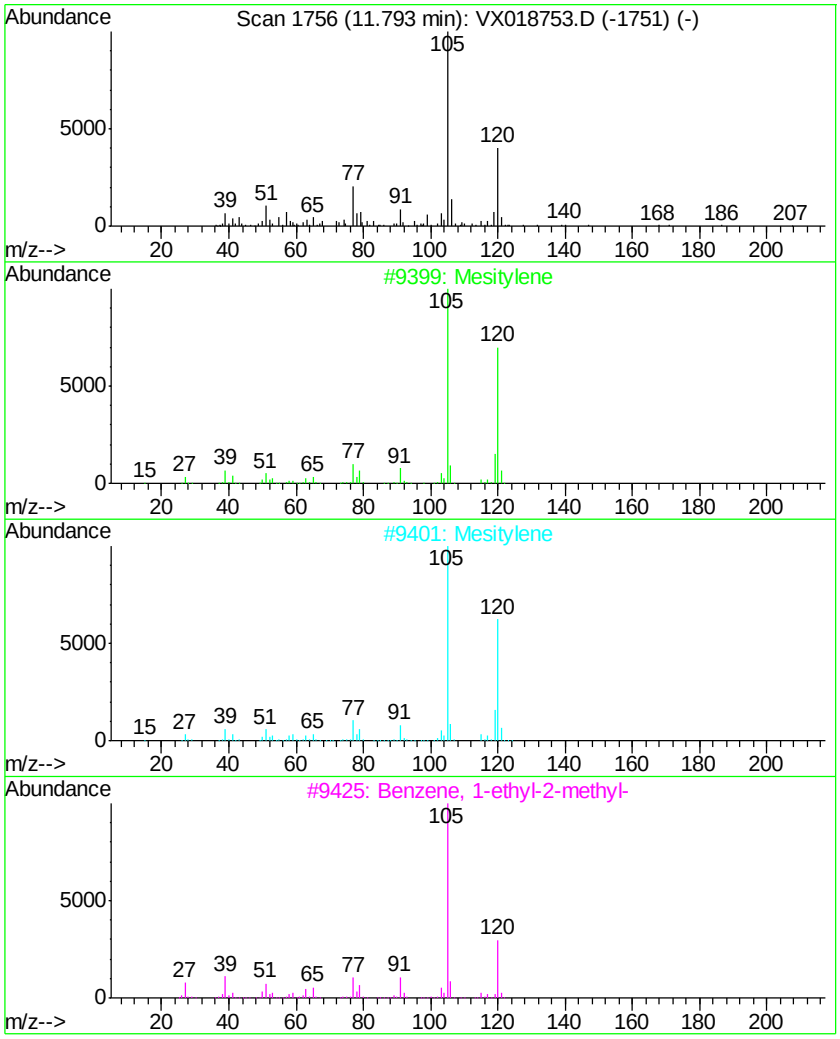
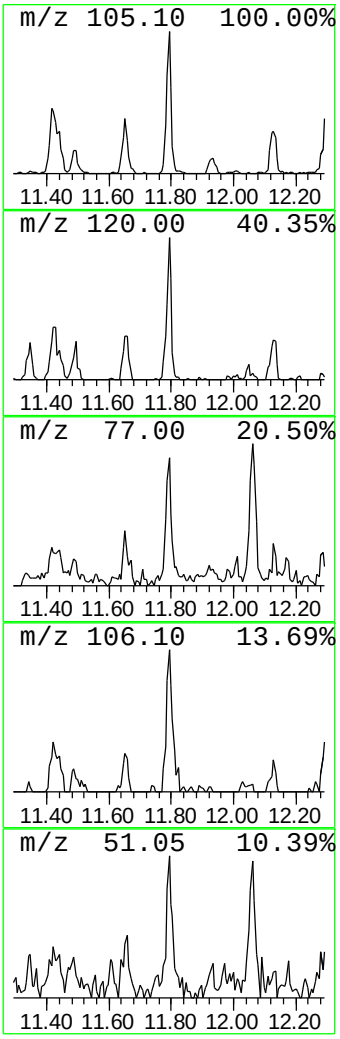
Peak Number 10 Mesitylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	0.69 ug/L	89407	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	81
2	Mesitylene	120	C9H12	000108-67-8	76
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	76
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	74
5	Mesitylene	120	C9H12	000108-67-8	72





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100220\
Data File : VX018753.D
Acq On : 02 Oct 2020 17:28
Operator : JC/SP
Sample : L4171-11
Misc : 25.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BG3Q1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

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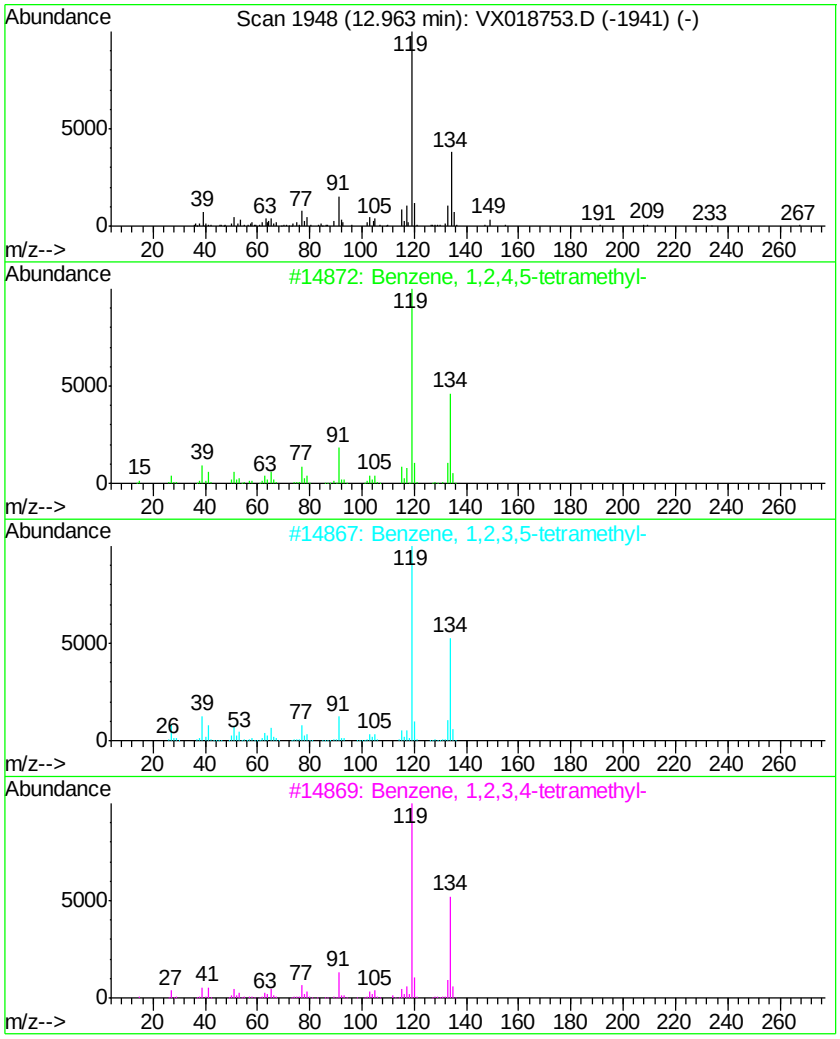
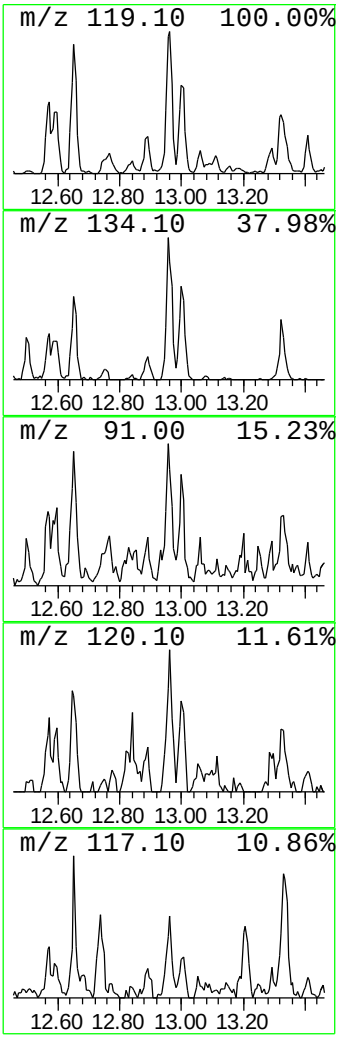
Peak Number 11 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	0.54 ug/L	70181	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	93
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	93
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	90
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100220\
Data File : VX018753.D
Acq On : 02 Oct 2020 17:28
Operator : JC/SP
Sample : L4171-11
Misc : 25.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BG3Q1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: Str...	1.16	0.6	ug/L	63167	1	6.83	508710	5.0
(DEL) Alkane: Str...	3.15	0.8	ug/L	85553	1	6.83	508710	5.0
(DEL) Alkane: Str...	3.50	1.2	ug/L	124863	1	6.83	508710	5.0
(DEL) Alkane: Cyc...	4.37	0.9	ug/L	94434	1	6.83	508710	5.0
2-Heptanone	10.81	0.6	ug/L	74635	2	10.10	627278	5.0
(DEL) Alkane: Cyc...	10.90	0.6	ug/L	76946	2	10.10	627278	5.0
Benzene, 1-ethyl-...	11.42	0.6	ug/L	70985	3	12.06	649197	5.0
4-Heptanone, 2,6-...	11.61	0.8	ug/L	104998	3	12.06	649197	5.0
Mesitylene	11.79	0.7	ug/L	89407	3	12.06	649197	5.0
Benzene, 1,2,4,5-...	12.96	0.5	ug/L	70181	3	12.06	649197	5.0