LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX050919\
Data File : VX009425.D
Acq On : 09 May 2019 14:14
Operator : JC/SP
Sample : K2748-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-8

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 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\82X042919W.M
Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.325	20	28	34	rBV2	48495	140802	3.97%	0.426%
2	1.404	38	41	46	rVB	49981	57979	1.63%	0.175%
3	1.788	100	104	122	rVB	96472	153776	4.33%	0.465%
4	2.001	134	139	146	rVB	29261	45027	1.27%	0.136%
5	2.605	229	238	251	rBV	24039	55030	1.55%	0.166%
6	2.843	261	277	282	rBV2	61686	168887	4.76%	0.511%
7	2.934	287	292	301	rVV	100673	218575	6.16%	0.661%
8	3.038	301	309	323	rVV	241685	542162	15.27%	1.640%
9	3.190	324	334	351	rVB	335997	831525	23.42%	2.515%
10	3.855	433	443	451	rBV2	14329	41782	1.18%	0.126%
11	4.403	524	533	546	rVB2	82431	249831	7.04%	0.756%
12	5.501	700	713	720	rBV	105728	312514	8.80%	0.945%
13	5.580	720	726	732	rVV2	33264	90847	2.56%	0.275%
14	5.659	732	739	749	rVV	169526	456096	12.85%	1.380%
15	5.940	771	785	795	rBV6	15953	55886	1.57%	0.169%
16	6.062	795	805	811	rVV	129808	339451	9.56%	1.027%
17	6.141	811	818	830	rVV	445017	1173770	33.06%	3.551%
18	6.342	831	851	862	rVV3	215553	700358	19.73%	2.119%
19	6.458	862	870	881	rVB7	17251	58639	1.65%	0.177%
20	6.860	924	936	946	rBV	304027	730608	20.58%	2.210%
21	7.458	1023	1034	1046	rBV4	50017	137828	3.88%	0.417%
22	8.055	1125	1132	1142	rVB	49312	98580	2.78%	0.298%
23	8.177	1144	1152	1162	rBV2	79901	205602	5.79%	0.622%
24	8.494	1198	1204	1210	rBV	57253	91911	2.59%	0.278%
25	8.567	1210	1216	1225	rVB	182249	290375	8.18%	0.878%
26	8.714	1227	1240	1246	rBV	655034	1038652	29.26%	3.142%
27	8.781	1246	1251	1259	rVB	902971	1409833	39.71%	4.265%
28	8.902	1264	1271	1280	rVB	185233	290135	8.17%	0.878%
29	9.500	1363	1369	1375	rBV2	39139	67548	1.90%	0.204%
30	9.622	1384	1389	1393	rBV	29739	41787	1.18%	0.126%
31	10.079	1457	1464	1466	rBV	135358	207781	5.85%	0.629%
32	10.110	1466	1469	1474	rVV	635979	820476	23.11%	2.482%
33	10.158	1474	1477	1480	rVV2	40029	55071	1.55%	0.167%
34	10.201	1480	1484	1487	rVV	179620	233932	6.59%	0.708%

LSC Area Percent Report

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ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-8

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 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\82X042919W.M
Title : SW846 8260

35	10.250	1487	1492	1498	rVV	2782188	3549911	100.00%	10.739%
36	10.299	1498	1500	1504	rVV2	31398	41203	1.16%	0.125%
37	10.360	1504	1510	1515	rVB	2265299	3081222	86.80%	9.321%
38	10.695	1560	1565	1575	rVB	1154342	1491771	42.02%	4.513%
39	10.841	1584	1589	1596	rBV2	64973	125440	3.53%	0.379%
40	10.951	1601	1607	1611	rVV3	24882	45934	1.29%	0.139%
41	11.018	1611	1618	1623	rVV	255575	338657	9.54%	1.024%
42	11.134	1631	1637	1642	rBV	515837	678299	19.11%	2.052%
43	11.225	1646	1652	1658	rVB2	86525	152821	4.30%	0.462%
44	11.323	1663	1668	1670	rVV	89979	121262	3.42%	0.367%
45	11.359	1670	1674	1681	rVV	591861	744274	20.97%	2.252%
46	11.433	1681	1686	1693	rVV	869194	1539243	43.36%	4.656%
47	11.506	1693	1698	1708	rVB	433718	543564	15.31%	1.644%
48	11.664	1714	1724	1732	rBV	729574	903608	25.45%	2.734%
49	11.804	1742	1747	1752	rVV	1819962	2130095	60.00%	6.444%
50	11.920	1760	1766	1767	rBV	72875	82596	2.33%	0.250%
51	11.939	1767	1769	1777	rVB2	84636	113771	3.20%	0.344%
52	12.024	1778	1783	1786	rBV	55170	65799	1.85%	0.199%
53	12.079	1786	1792	1797	rVB	608985	786513	22.16%	2.379%
54	12.140	1797	1802	1809	rBV	499641	617123	17.38%	1.867%
55	12.298	1820	1828	1835	rBV2	1029454	1483993	41.80%	4.489%
56	12.371	1835	1840	1848	rVB3	213599	354371	9.98%	1.072%
57	12.457	1848	1854	1860	rBV2	45142	67700	1.91%	0.205%
58	12.518	1860	1864	1870	rVB	111903	133790	3.77%	0.405%
59	12.585	1870	1875	1877	rBV	149300	185533	5.23%	0.561%
60	12.609	1877	1879	1884	rVV	91517	96366	2.71%	0.292%
61	12.664	1884	1888	1892	rVV	177105	222103	6.26%	0.672%
62	12.701	1892	1894	1899	rVV	48730	53136	1.50%	0.161%
63	12.755	1899	1903	1910	rVB2	134528	206619	5.82%	0.625%
64	12.902	1922	1927	1932	rVB	69652	93913	2.65%	0.284%
65	12.975	1932	1939	1942	rBV	162231	203553	5.73%	0.616%
66	13.018	1942	1946	1951	rVB	222001	268677	7.57%	0.813%
67	13.079	1951	1956	1962	rBV7	16153	37589	1.06%	0.114%
68	13.225	1975	1980	1984	rVB	115450	134254	3.78%	0.406%
69	13.347	1995	2000	2006	rVB2	262954	341137	9.61%	1.032%
70	13.475	2017	2021	2028	rVB	34821	48577	1.37%	0.147%
71	13.828	2073	2079	2085	rBV	366112	462243	13.02%	1.398%

LSC Area Percent Report

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ALS Vial : 11 Sample Multiplier: 1

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ClientSampleId : MW-8

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 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\82X042919W.M
Title : SW846 8260

72	14.694	2216	2221	2226	rVB	36234	43772	1.23%	0.132%
73	14.834	2237	2244	2249	rBV	39232	52943	1.49%	0.160%

Sum of corrected areas: 33056431

LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX050919\
Data File : VX009425.D
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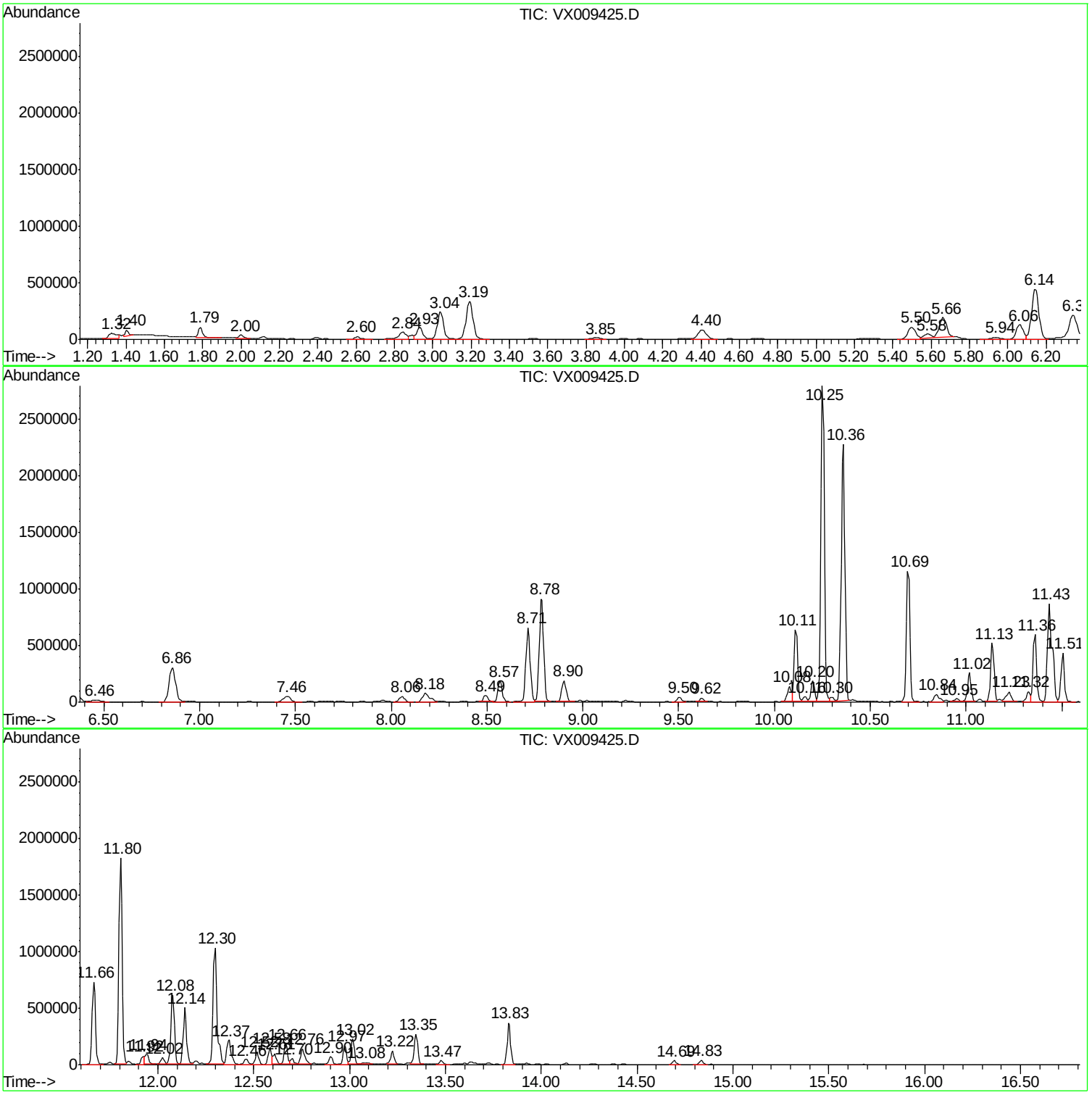
Operator : JC/SP
Sample : K2748-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX050919\
Data File : VX009425.D
Acq On : 09 May 2019 14:14
Operator : JC/SP
Sample : K2748-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

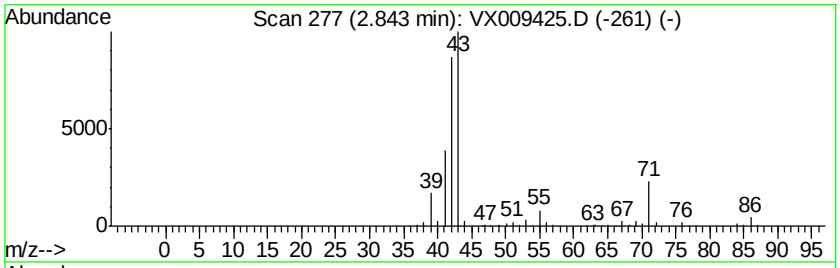
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

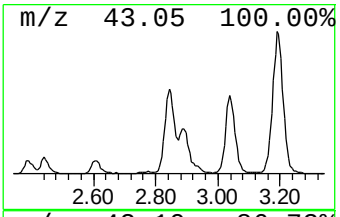
****

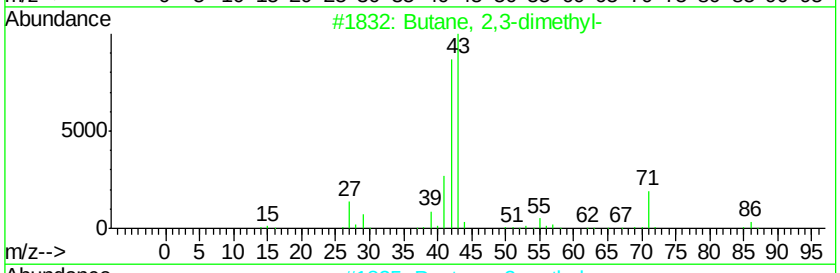
Peak Number 1 Butane, 2,3-dimethyl- Concentration Rank 9

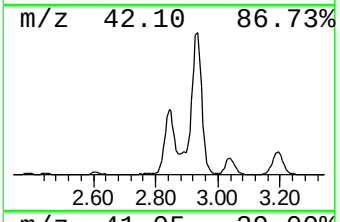
R.T.	EstConc	Area	Relative to ISTD	R.T.
2.84	18.51 ug/l	168887	Pentafluorobenzene	5.67

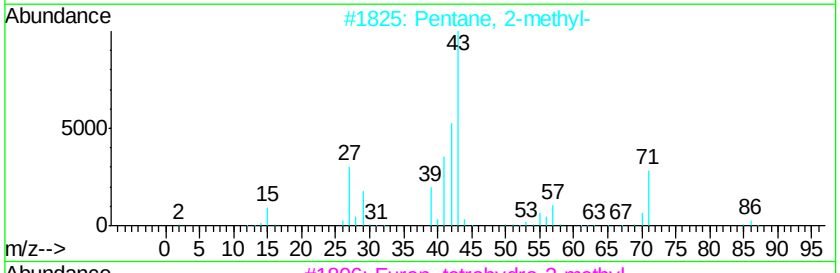
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	47
3	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	14
4	Oxalic acid, pentyl propyl ester	202	C10H18O4	1000309-25-7	9
5	1-Hexene, 4,5-dimethyl-	112	C8H16	016106-59-5	9

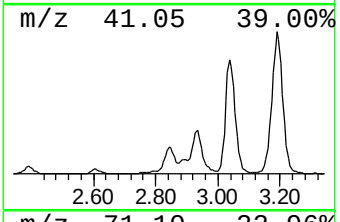


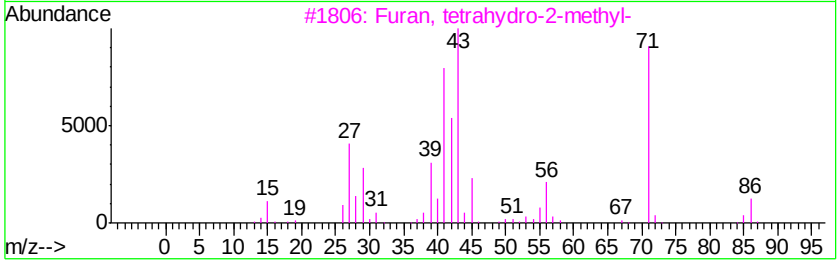


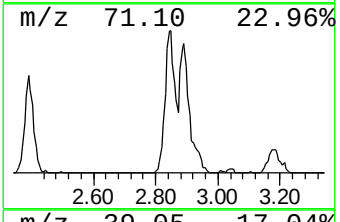


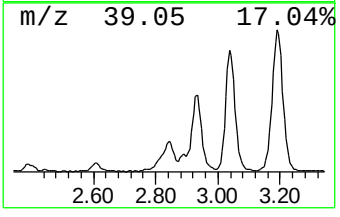












Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX050919\
Data File : VX009425.D
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Operator : JC/SP
Sample : K2748-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

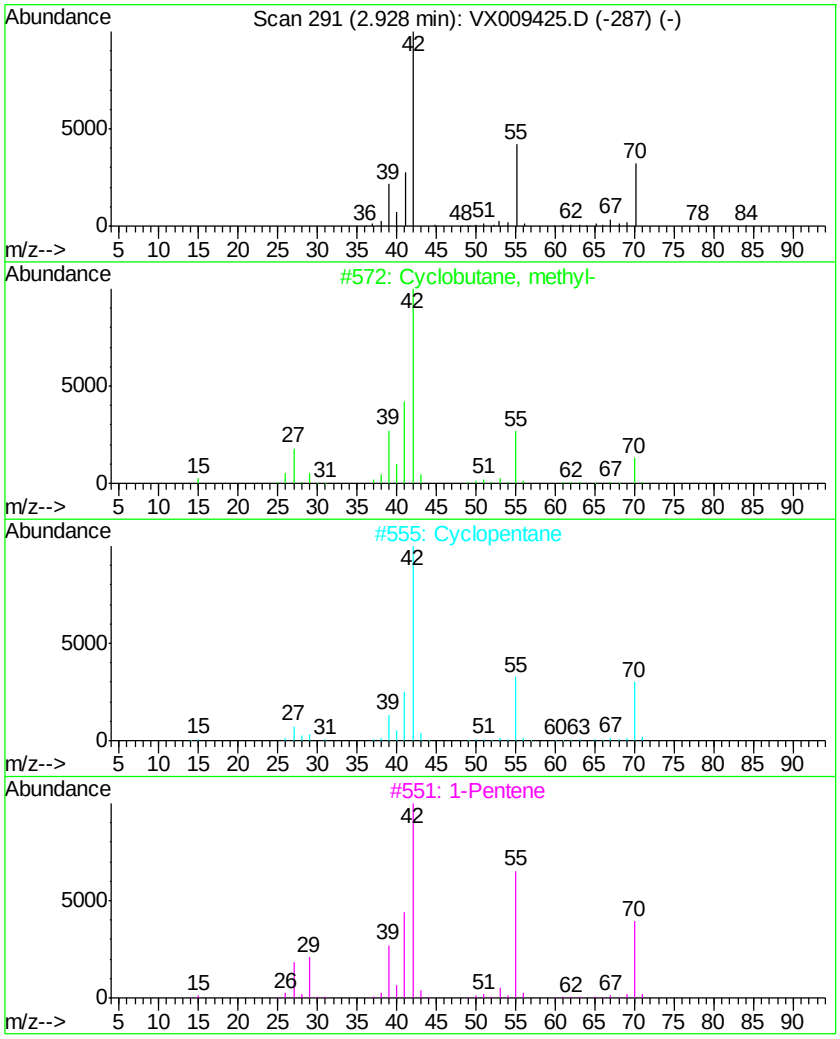
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

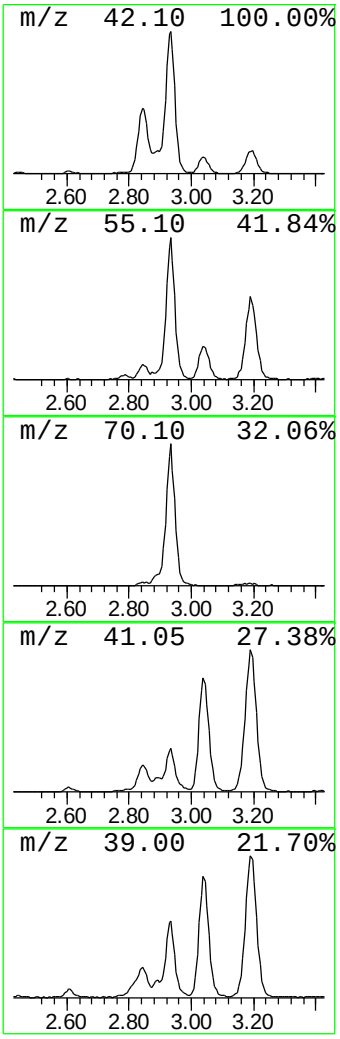
****

Peak Number 2 Cyclobutane, methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.93	23.96 ug/l	218575	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclobutane, methyl-	70	C5H10	000598-61-8	86
2	Cyclopentane	70	C5H10	000287-92-3	80
3	1-Pentene	70	C5H10	000109-67-1	46
4	Cyclopropane, ethyl-	70	C5H10	001191-96-4	23
5	Methyl propargyl ether	70	C4H6O	000627-41-8	7





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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
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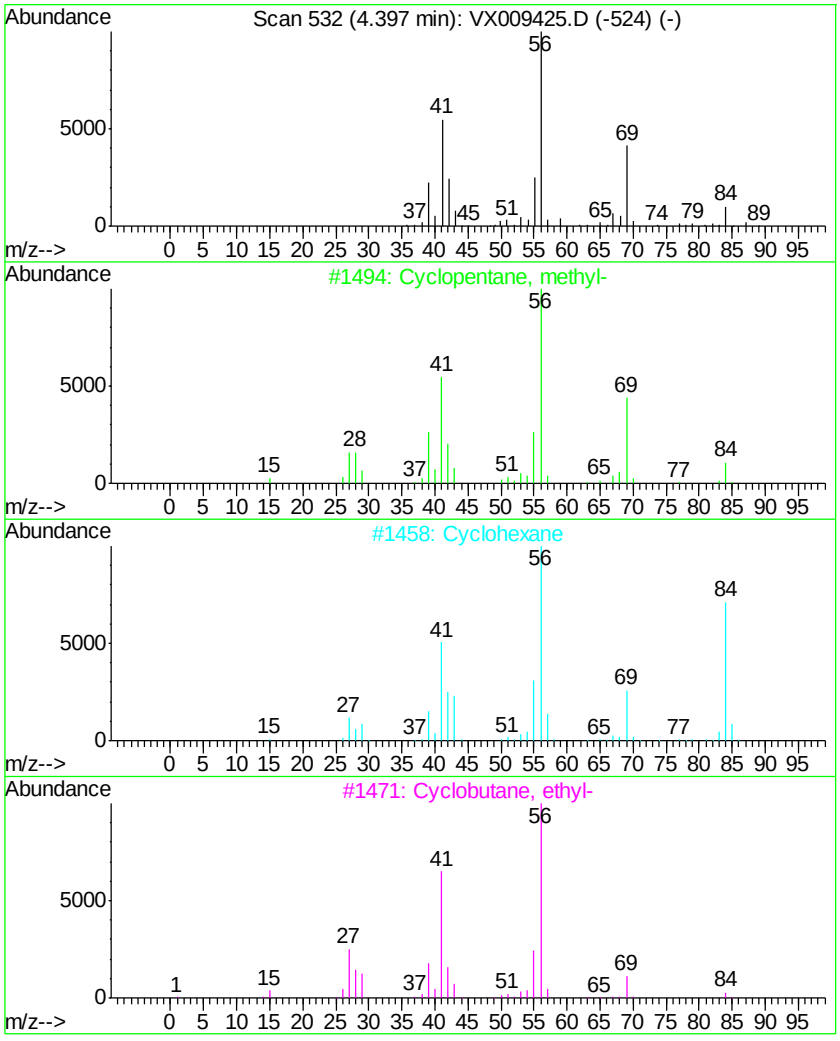
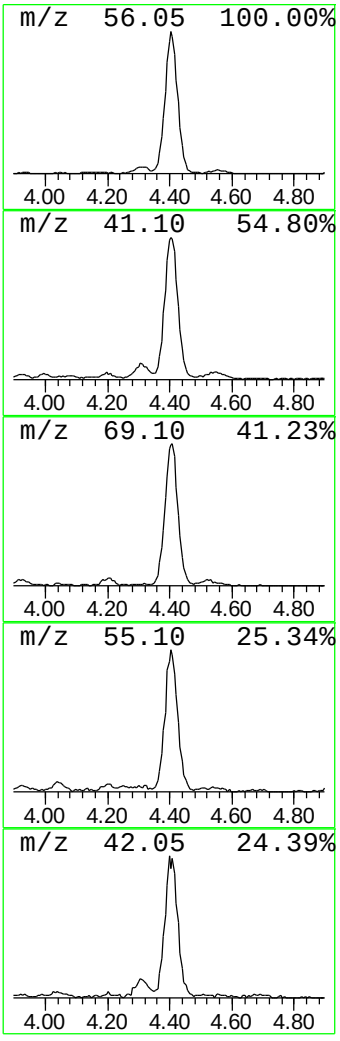
****

Peak Number 3 Cyclopentane, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	27.39 ug/l	249831	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	93
2	Cyclohexane	84	C6H12	000110-82-7	72
3	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	50
5	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	40





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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

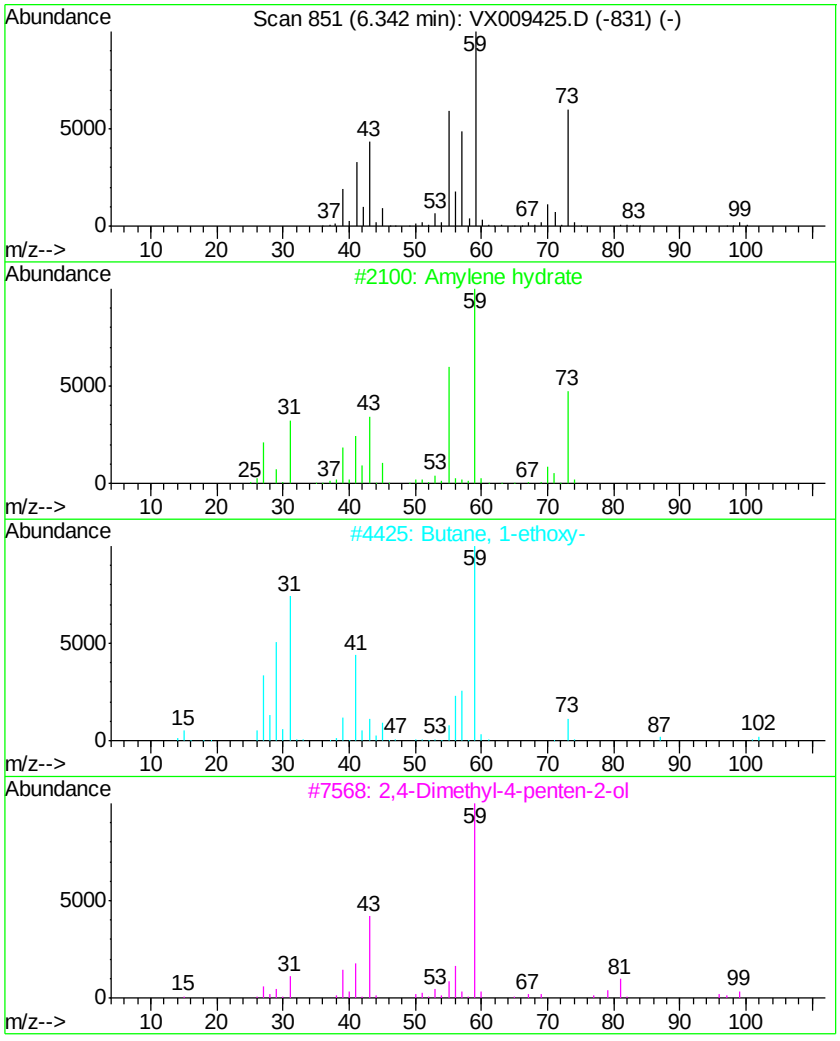
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

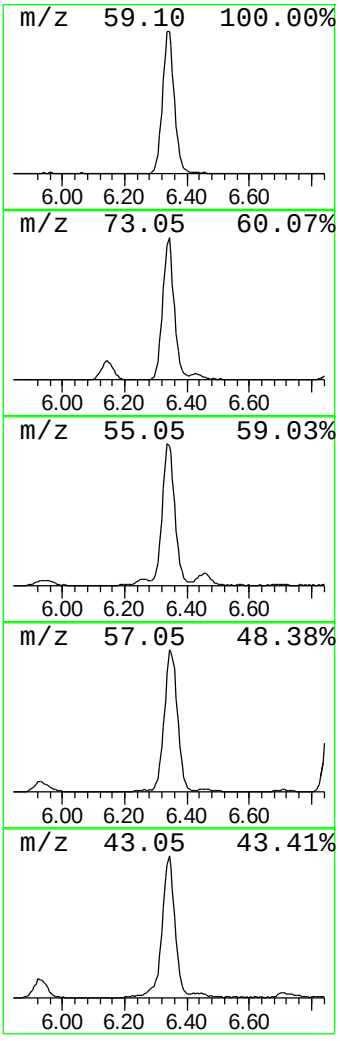
****

Peak Number 4 Amylene hydrate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.34	47.93 ug/l	700358	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Amylene hydrate	88	C5H12O	000075-85-4	72
2	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	39
3	2,4-Dimethyl-4-penten-2-ol	114	C7H14O	019781-53-4	38
4	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	38
5	3-Hexanol	102	C6H14O	000623-37-0	38





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ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

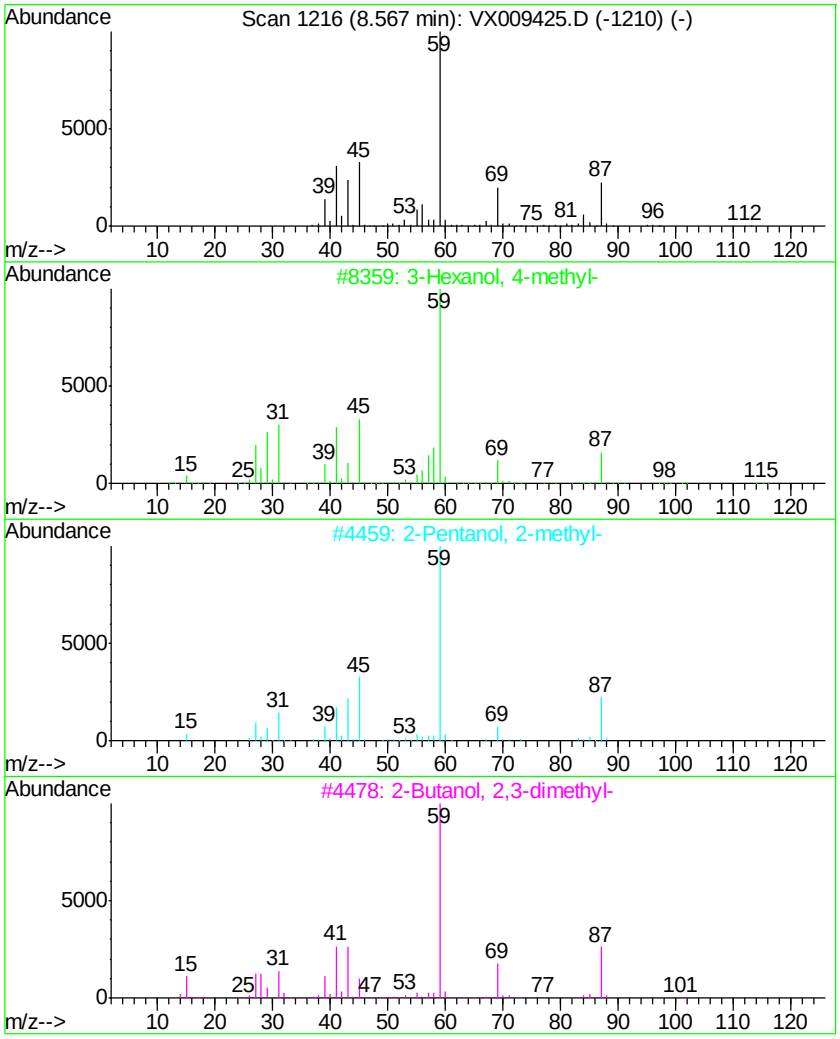
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

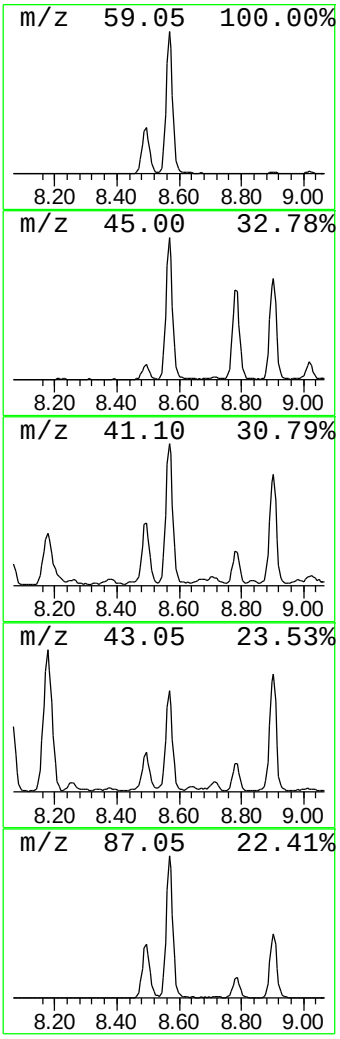
****

Peak Number 5 3-Hexanol, 4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.57	17.70 ug/l	290375	Chlorobenzene-d5	10.11

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	78
2	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	78
3	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	50
4	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	45
5	Propanoic acid, 3-nitro-, methyl...	133	C4H7NO4	020497-95-4	39





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX050919\
Data File : VX009425.D
Acq On : 09 May 2019 14:14
Operator : JC/SP
Sample : K2748-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

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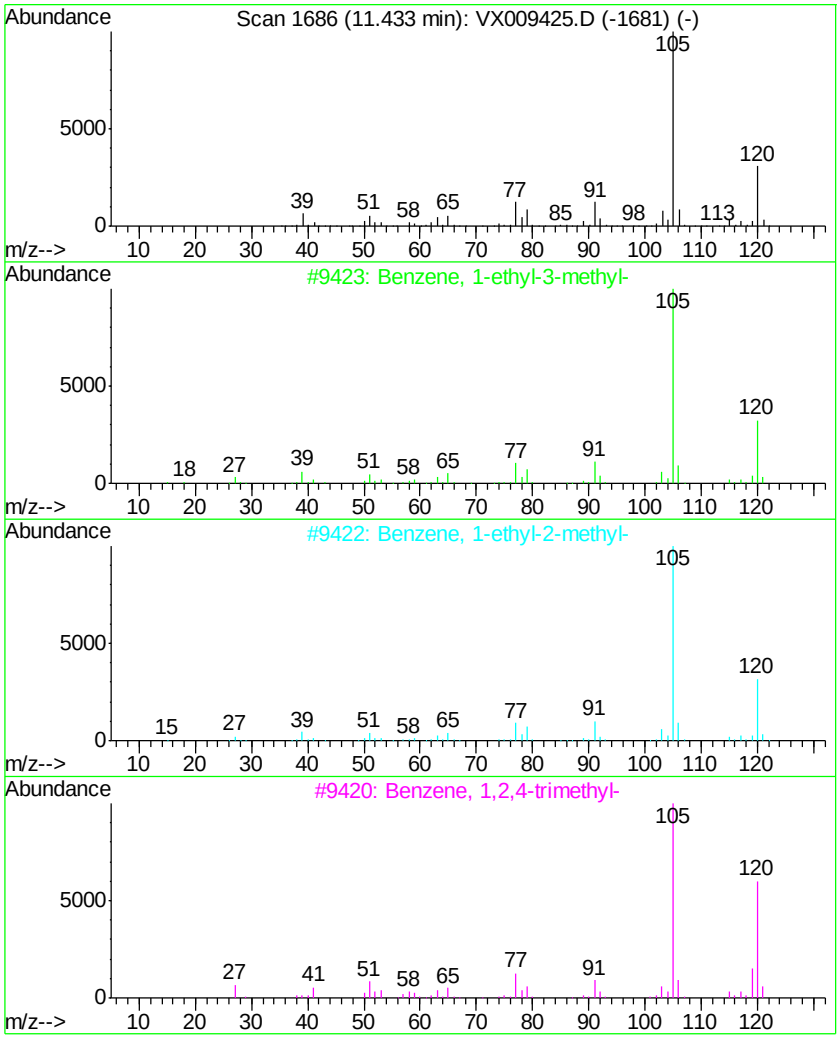
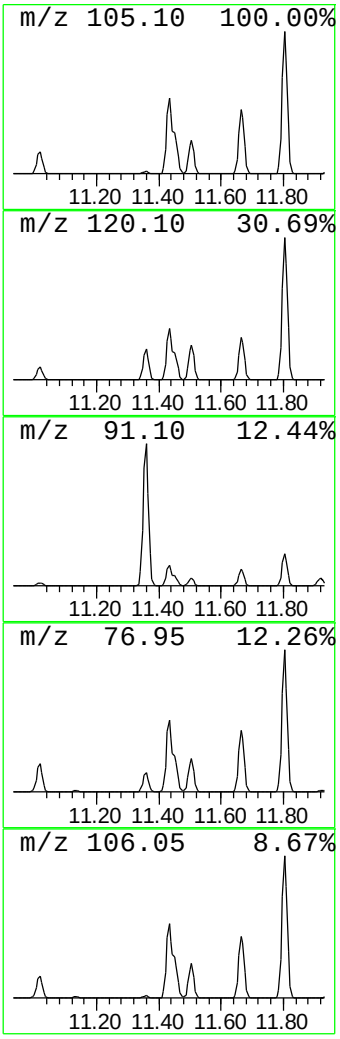
Peak Number 6 Benzene, 1-ethyl-3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	97.85 ug/l	1539240	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX050919\
Data File : VX009425.D
Acq On : 09 May 2019 14:14
Operator : JC/SP
Sample : K2748-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

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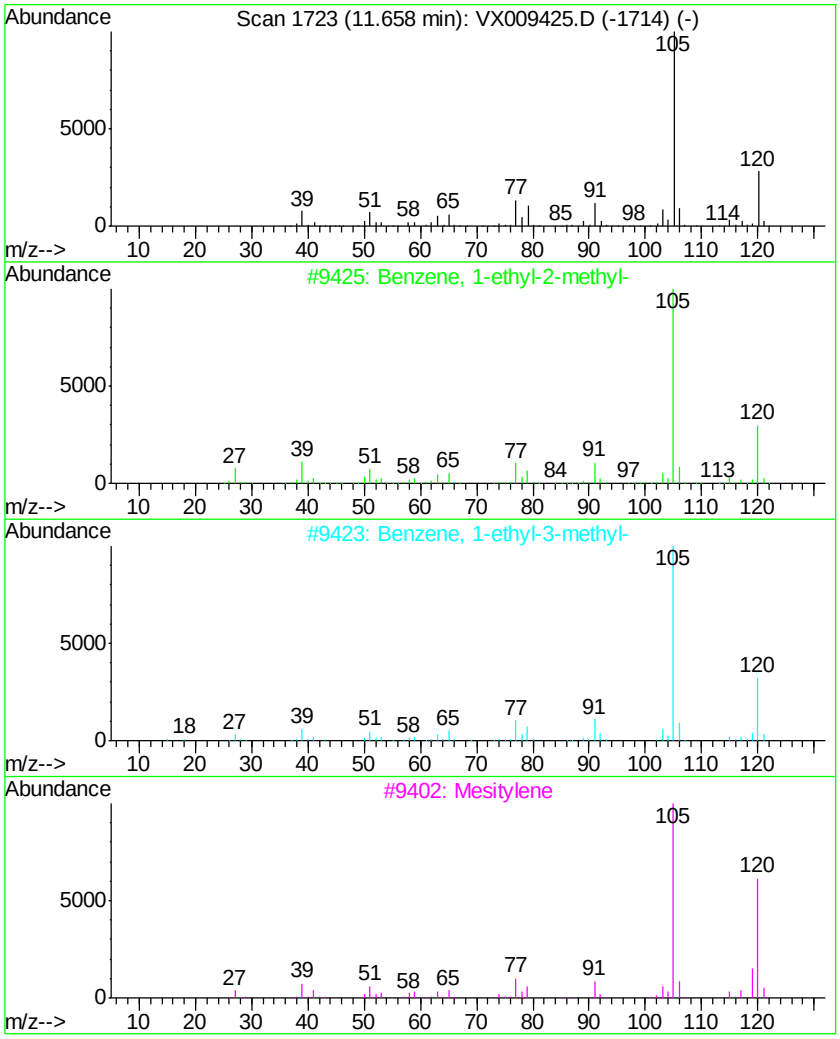
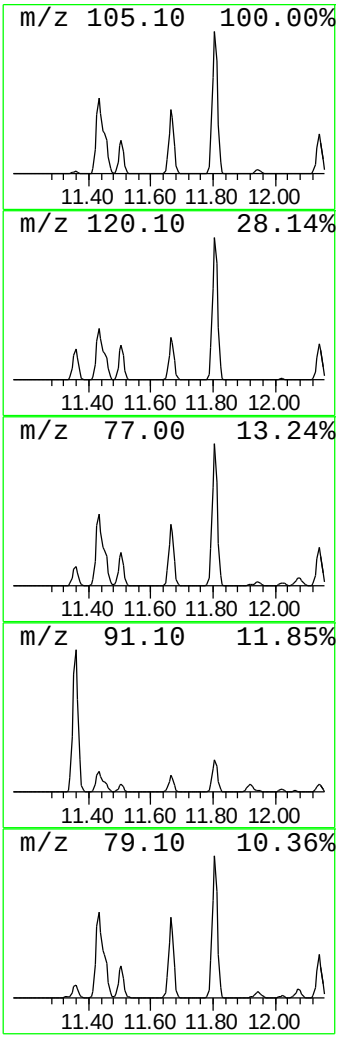
Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	57.44 ug/l	903608	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX050919\
Data File : VX009425.D
Acq On : 09 May 2019 14:14
Operator : JC/SP
Sample : K2748-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

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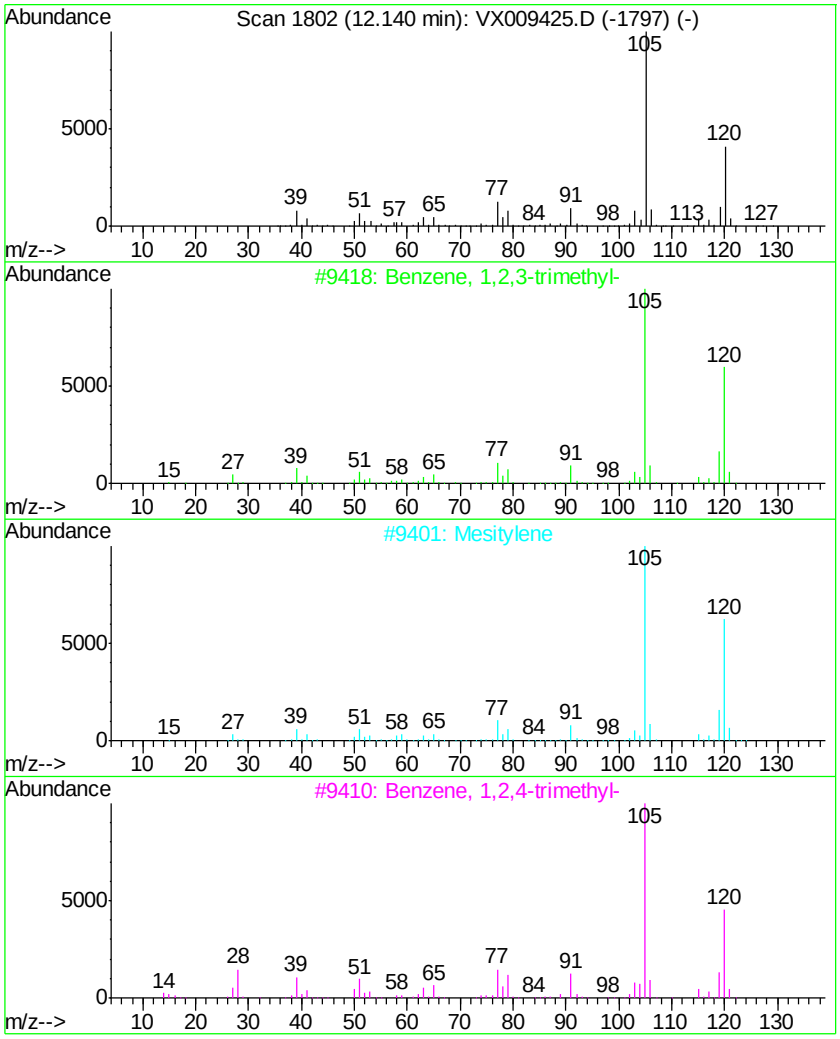
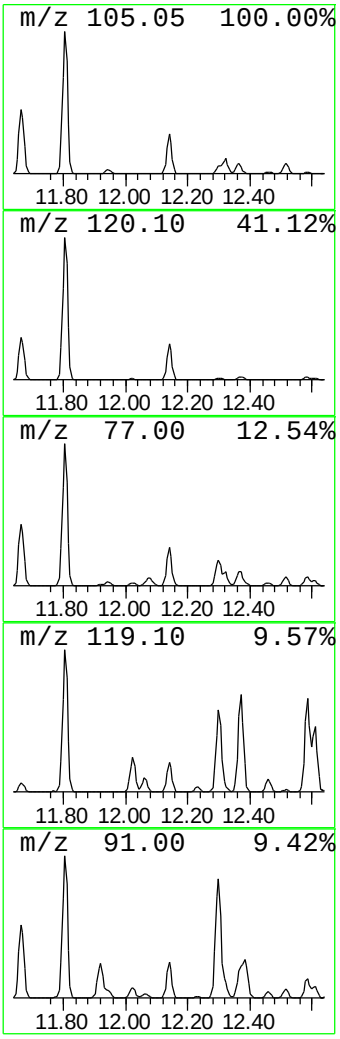
Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	39.23 ug/l	617123	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Mesitylene	120	C9H12	000108-67-8	94
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX050919\
Data File : VX009425.D
Acq On : 09 May 2019 14:14
Operator : JC/SP
Sample : K2748-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

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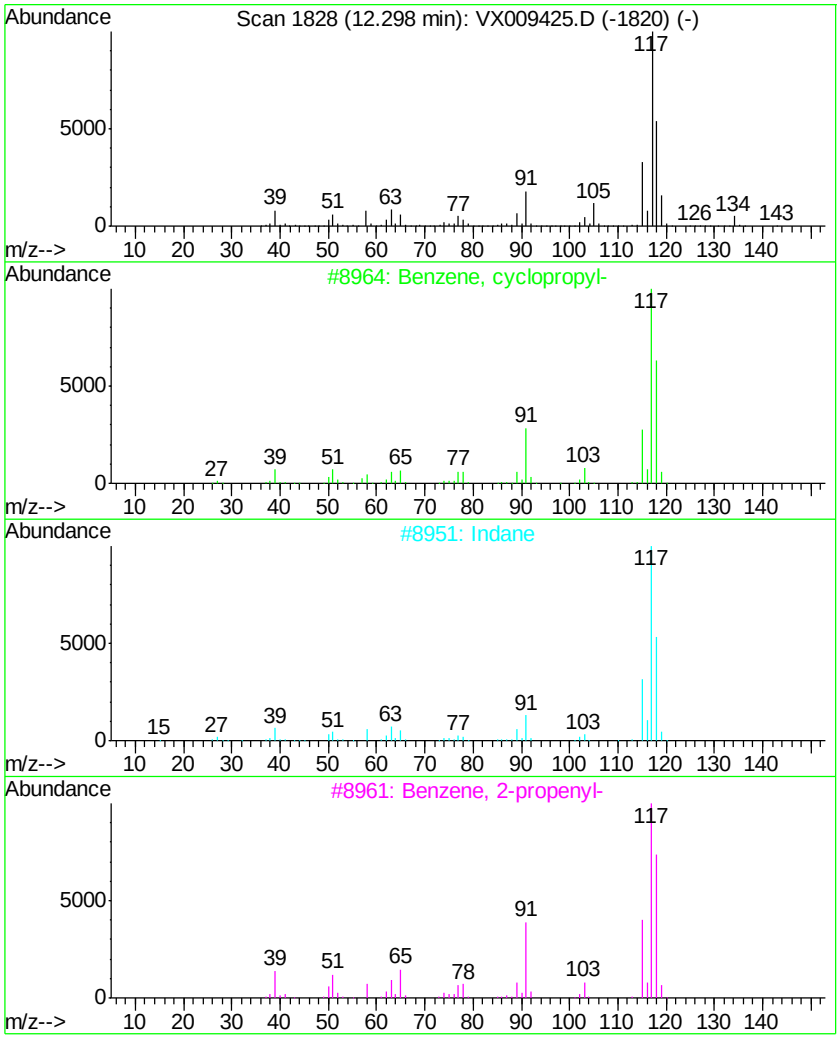
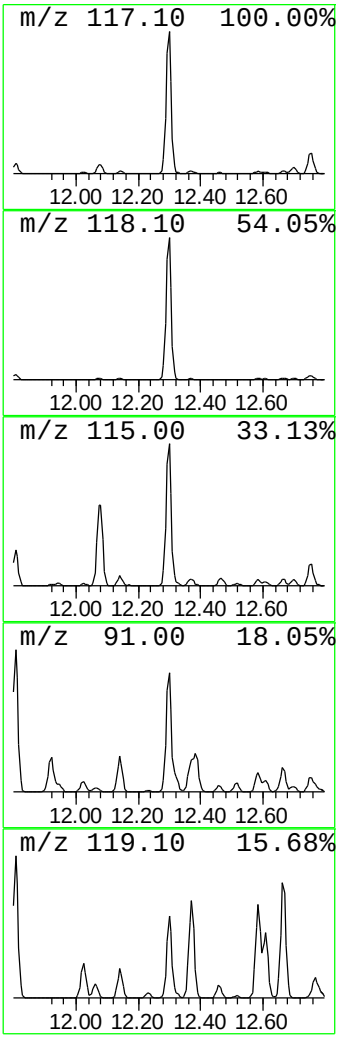
Peak Number 9 Benzene, cyclopropyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	94.34 ug/l	1483990	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
2	Indane	118	C9H10	000496-11-7	76
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
4	Deltacyclene	118	C9H10	007785-10-6	58
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	55





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX050919\
Data File : VX009425.D
Acq On : 09 May 2019 14:14
Operator : JC/SP
Sample : K2748-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

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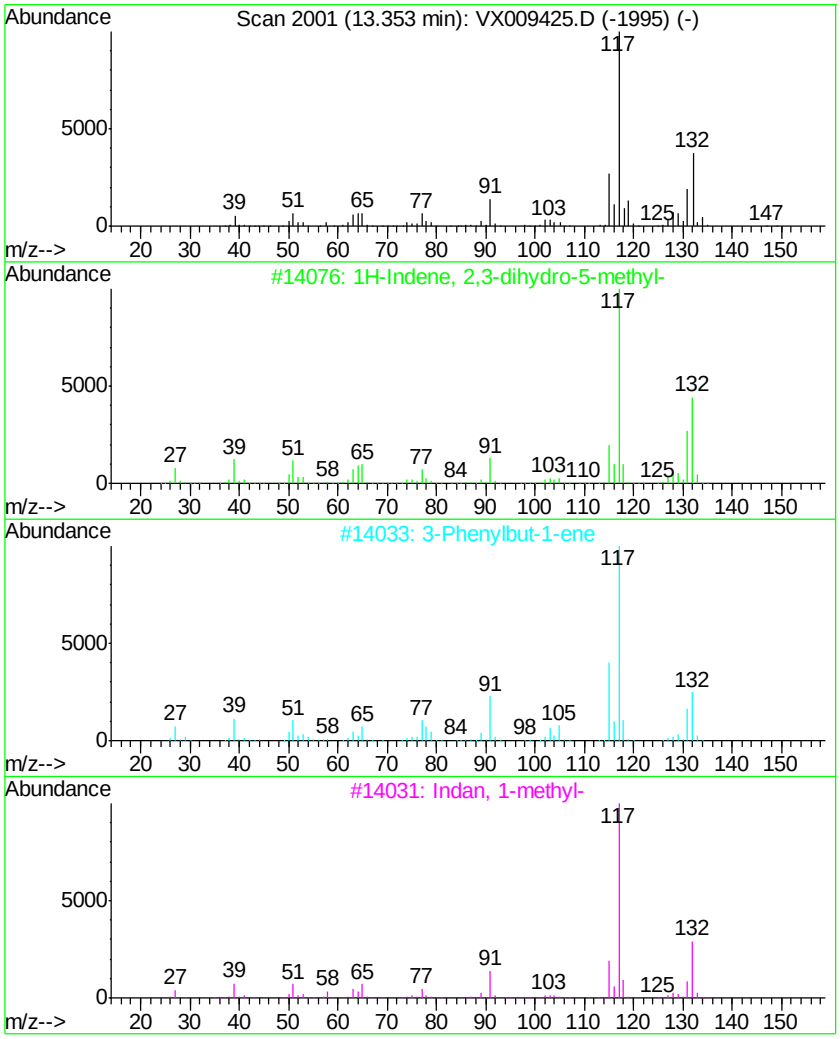
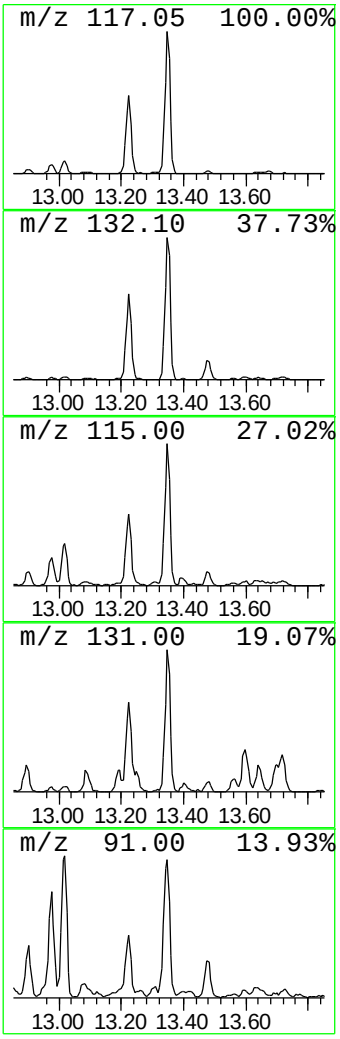
Peak Number 10 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	21.69 ug/l	341137	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3	Indan, 1-methyl-	132	C10H12	000767-58-8	76
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX050919\
Data File : VX009425.D
Acq On : 09 May 2019 14:14
Operator : JC/SP
Sample : K2748-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2,3-dimet...	2.84	18.5	ug/l	168887	1	5.67	456096	50.0
Cyclobutane, methyl-	2.93	24.0	ug/l	218575	1	5.67	456096	50.0
Cyclopentane, met...	4.40	27.4	ug/l	249831	1	5.67	456096	50.0
Amylene hydrate	6.34	47.9	ug/l	700358	2	6.86	730608	50.0
3-Hexanol, 4-methyl-	8.57	17.7	ug/l	290375	3	10.11	820476	50.0
Benzene, 1-ethyl-...	11.43	97.8	ug/l	1539240	4	12.08	786513	50.0
Benzene, 1-ethyl-...	11.66	57.4	ug/l	903608	4	12.08	786513	50.0
Benzene, 1,2,3-tr...	12.14	39.2	ug/l	617123	4	12.08	786513	50.0
Benzene, cyclopro...	12.30	94.3	ug/l	1483990	4	12.08	786513	50.0
1H-Indene, 2,3-di...	13.35	21.7	ug/l	341137	4	12.08	786513	50.0