LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010726.D
Acq On : 09 Jul 2019 03:53
Operator : JC/SP
Sample : K3690-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW082-190703

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\82X061919W.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	1.276	16	20	26	rBV	49306	65154	4.73%	0.548%
2	1.404	38	41	47	rBV	24729	28050	2.03%	0.236%
3	1.788	101	104	112	rVB	16125	24085	1.75%	0.202%
4	2.843	270	277	289	rBV	32958	75837	5.50%	0.637%
5	3.178	323	332	338	rBV3	7772	18073	1.31%	0.152%
6	4.306	508	517	526	rBV3	6754	19528	1.42%	0.164%
7	4.598	550	565	579	rBV	173096	491510	35.66%	4.131%
8	5.501	702	713	724	rBV	128527	371135	26.92%	3.120%
9	5.659	729	739	761	rVB	221575	650519	47.19%	5.468%
10	6.062	795	805	815	rBV	125582	329807	23.93%	2.772%
11	6.348	843	852	863	rVB3	35766	107623	7.81%	0.905%
12	6.860	926	936	950	rBV	334137	779923	56.58%	6.556%
13	7.220	988	995	1001	rBV3	8700	20667	1.50%	0.174%
14	7.452	1026	1033	1041	rVB9	7319	18920	1.37%	0.159%
15	7.683	1064	1071	1078	rVB2	8102	15533	1.13%	0.131%
16	8.055	1121	1132	1138	rBV	24411	51273	3.72%	0.431%
17	8.177	1145	1152	1161	rVB2	77677	154461	11.21%	1.298%
18	8.713	1229	1240	1249	rBV	706134	1096188	79.52%	9.214%
19	9.335	1334	1342	1348	rBV	21732	34539	2.51%	0.290%
20	10.110	1463	1469	1471	rBV	701649	993035	72.04%	8.347%
21	10.134	1471	1473	1479	rVV	1083077	1378468	100.00%	11.587%
22	10.183	1479	1481	1486	rVB	10545	14198	1.03%	0.119%
23	10.335	1497	1506	1516	rVV4	25256	48605	3.53%	0.409%
24	10.439	1516	1523	1530	rVB10	5948	13835	1.00%	0.116%
25	10.512	1530	1535	1541	rBV2	15444	23307	1.69%	0.196%
26	10.622	1546	1553	1555	rBV4	11103	19967	1.45%	0.168%
27	10.658	1555	1559	1564	rVB4	39557	60270	4.37%	0.507%
28	10.707	1564	1567	1572	rVB3	10349	14746	1.07%	0.124%
29	10.792	1576	1581	1585	rBV3	13678	27504	2.00%	0.231%
30	10.835	1585	1588	1594	rVB	19644	26489	1.92%	0.223%
31	10.914	1596	1601	1604	rBV3	17309	30885	2.24%	0.260%
32	10.969	1608	1610	1614	rVB	12141	14720	1.07%	0.124%
33	11.018	1614	1618	1620	rBV	14762	20250	1.47%	0.170%
34	11.061	1620	1625	1632	rVV6	24727	50413	3.66%	0.424%

LSC Area Percent Report

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Operator : JC/SP
Sample : K3690-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW082-190703

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\82X061919W.M
Title : SW846 8260

35	11.134	1632	1637	1642	rVV	596546	747065	54.20%	6.279%
36	11.183	1642	1645	1648	rVV2	51667	69354	5.03%	0.583%
37	11.244	1652	1655	1658	rVV2	20840	28964	2.10%	0.243%
38	11.280	1658	1661	1664	rVB	20535	24124	1.75%	0.203%
39	11.390	1675	1679	1684	rBV3	12891	18543	1.35%	0.156%
40	11.445	1684	1688	1692	rBV2	27217	39488	2.86%	0.332%
41	11.500	1692	1697	1700	rVV5	12810	17318	1.26%	0.146%
42	11.579	1706	1710	1716	rVB7	16002	30084	2.18%	0.253%
43	11.670	1718	1725	1731	rBV5	33110	60861	4.42%	0.512%
44	11.768	1731	1741	1744	rBV	106120	152236	11.04%	1.280%
45	11.804	1744	1747	1751	rVV3	18539	27492	1.99%	0.231%
46	11.914	1758	1765	1768	rBV	125322	196387	14.25%	1.651%
47	11.945	1768	1770	1775	rVB	71815	78072	5.66%	0.656%
48	12.024	1777	1783	1786	rVB4	11318	19191	1.39%	0.161%
49	12.079	1786	1792	1798	rBV	704421	943242	68.43%	7.928%
50	12.170	1803	1807	1812	rVV2	15507	27579	2.00%	0.232%
51	12.231	1812	1817	1823	rVB	344508	410955	29.81%	3.454%
52	12.298	1823	1828	1833	rBV	91306	115232	8.36%	0.969%
53	12.371	1835	1840	1848	rVB5	33268	68554	4.97%	0.576%
54	12.469	1848	1856	1861	rBV4	13072	27619	2.00%	0.232%
55	12.658	1878	1887	1890	rBV4	26143	54497	3.95%	0.458%
56	12.713	1890	1896	1899	rVV2	86343	166756	12.10%	1.402%
57	12.755	1899	1903	1909	rVV2	202288	345427	25.06%	2.904%
58	12.816	1909	1913	1917	rVV	67670	89646	6.50%	0.754%
59	12.871	1917	1922	1923	rVV	55297	75712	5.49%	0.636%
60	12.896	1923	1926	1931	rVV	191259	231857	16.82%	1.949%
61	12.957	1931	1936	1938	rVV	56928	91753	6.66%	0.771%
62	12.975	1938	1939	1945	rVB2	53107	49896	3.62%	0.419%
63	13.085	1952	1957	1962	rVB	65471	90742	6.58%	0.763%
64	13.188	1970	1974	1978	rBV	43845	53293	3.87%	0.448%
65	13.249	1980	1984	1988	rVV	69692	86159	6.25%	0.724%
66	13.304	1988	1993	1996	rVV	19863	30777	2.23%	0.259%
67	13.347	1996	2000	2005	rVV2	61736	83273	6.04%	0.700%
68	13.402	2005	2009	2014	rVV2	56146	74661	5.42%	0.628%
69	13.481	2014	2022	2027	rVB4	25248	53086	3.85%	0.446%
70	13.597	2036	2041	2045	rVB3	16221	25697	1.86%	0.216%
71	13.639	2045	2048	2051	rBV3	15416	19998	1.45%	0.168%

LSC Area Percent Report

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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW082-190703

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

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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Title : SW846 8260

72	13.670	2051	2053	2055	rBV2	13241	15498	1.12%	0.130%
73	13.853	2077	2083	2088	rVB6	7630	16794	1.22%	0.141%
74	13.908	2088	2092	2096	rBV4	10345	14841	1.08%	0.125%
75	13.981	2101	2104	2108	rVV3	12254	16184	1.17%	0.136%
76	14.840	2240	2245	2251	rVB3	9307	18478	1.34%	0.155%

Sum of corrected areas: 11896902

LSC Report - Integrated Chromatogram

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Data File : VX010726.D
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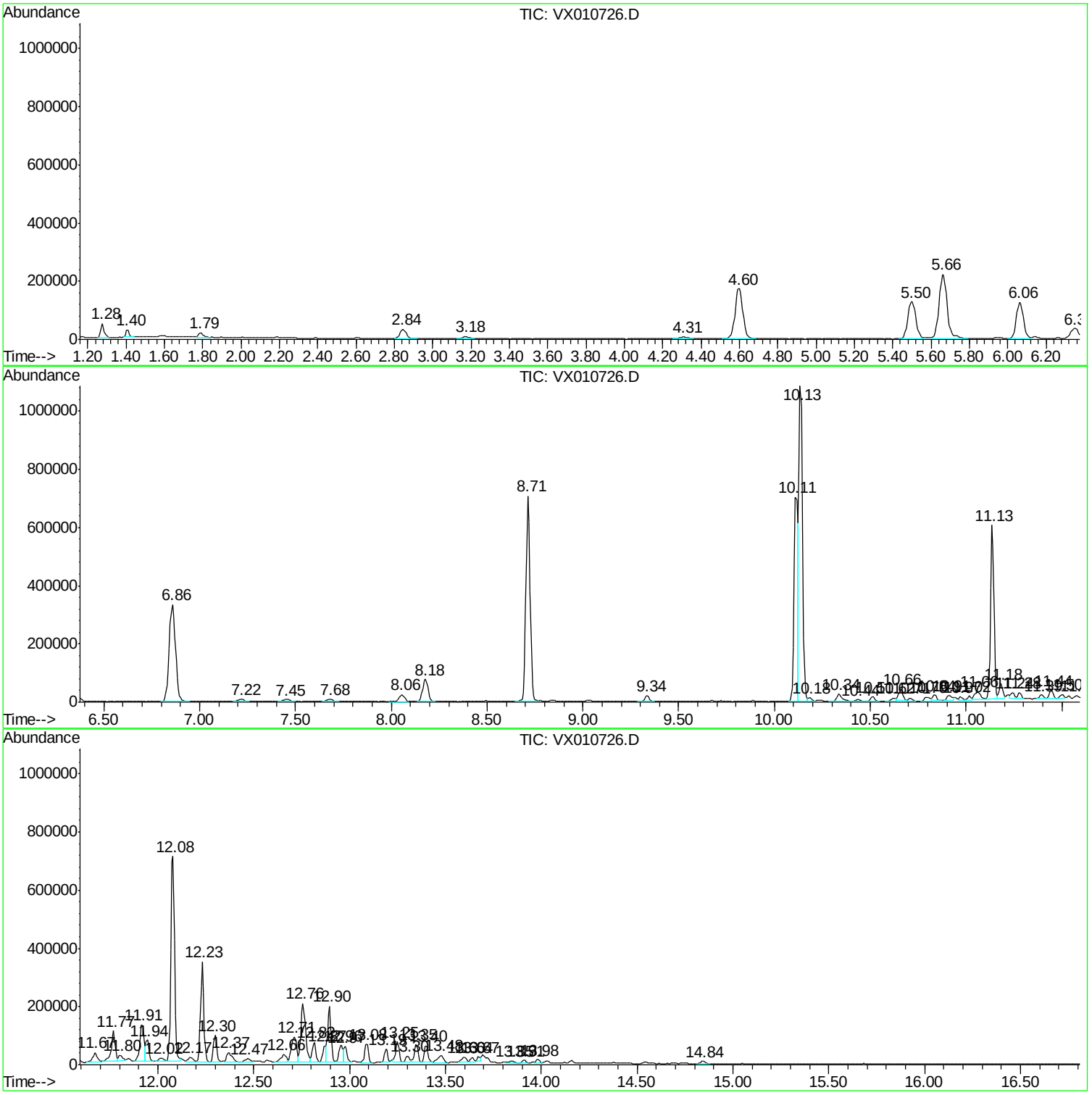
Operator : JC/SP
Sample : K3690-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SS038RW082-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010726.D
Acq On : 09 Jul 2019 03:53
Operator : JC/SP
Sample : K3690-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW082-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

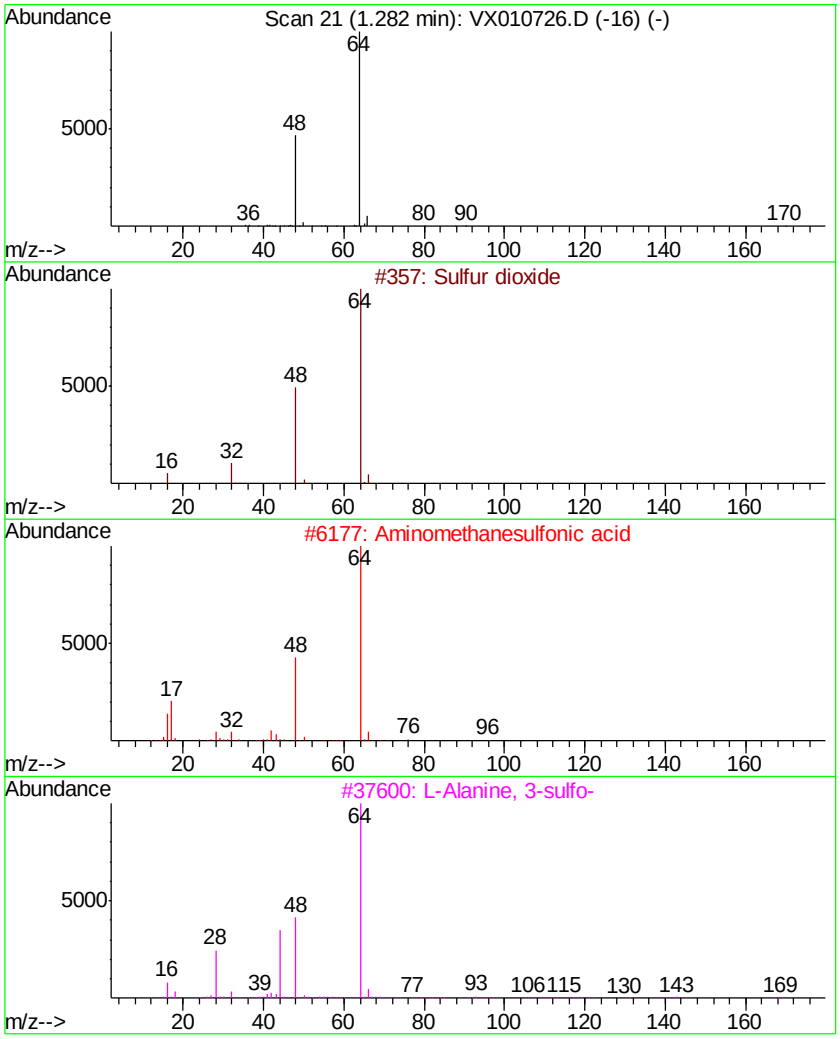
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

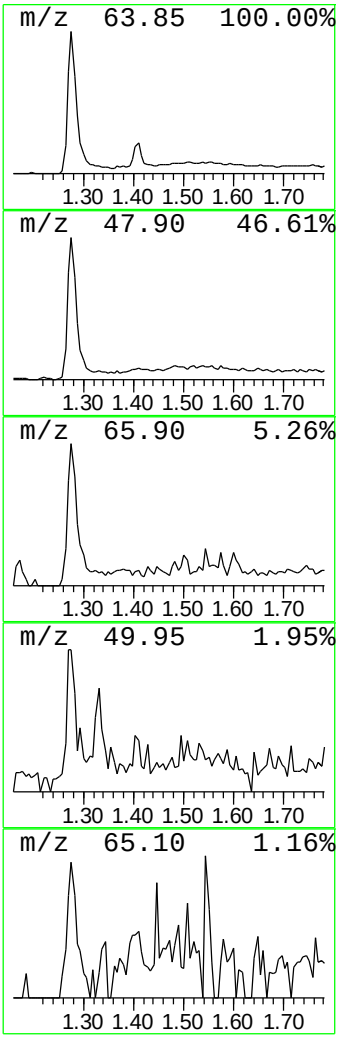
****

Peak Number 1 Sulfur dioxide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.28	5.01 ug/l	65154	Pentafluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfur dioxide	64	O2S	007446-09-5	90
2	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	64
3	L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
4	Ethyl Chloride	64	C2H5Cl	000075-00-3	3
5	Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	3





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
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Acq On : 09 Jul 2019 03:53
Operator : JC/SP
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW082-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

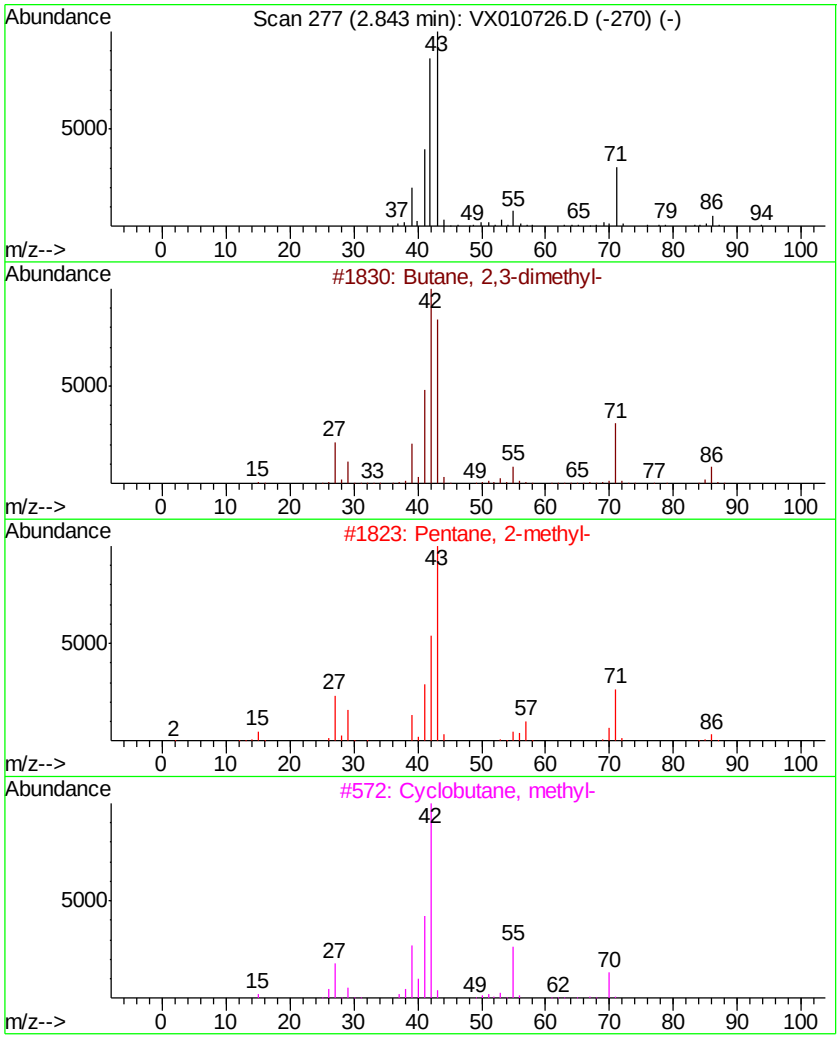
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

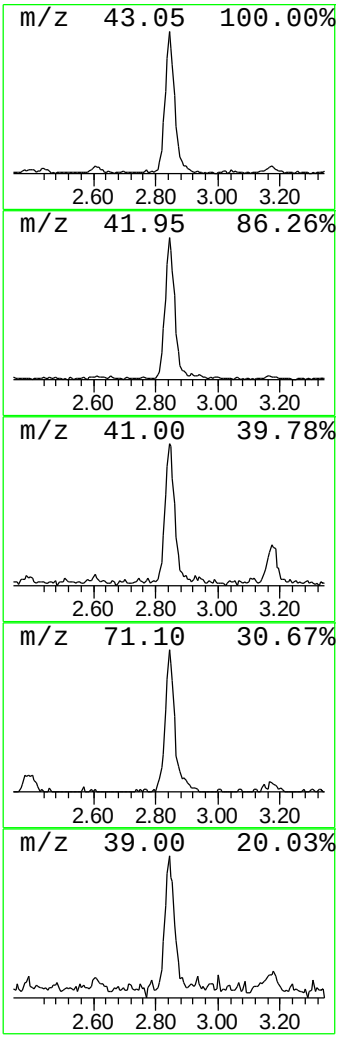
****

Peak Number 2 Butane, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.84	5.83 ug/l	75837	Pentafluorobenzene	5.66

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	50
3	Cyclobutane, methyl-	70	C5H10	000598-61-8	50
4	Tetrahydrofuran	72	C4H8O	000109-99-9	38
5	Butane, 2-methyl-	72	C5H12	000078-78-4	36





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010726.D
Acq On : 09 Jul 2019 03:53
Operator : JC/SP
Sample : K3690-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW082-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

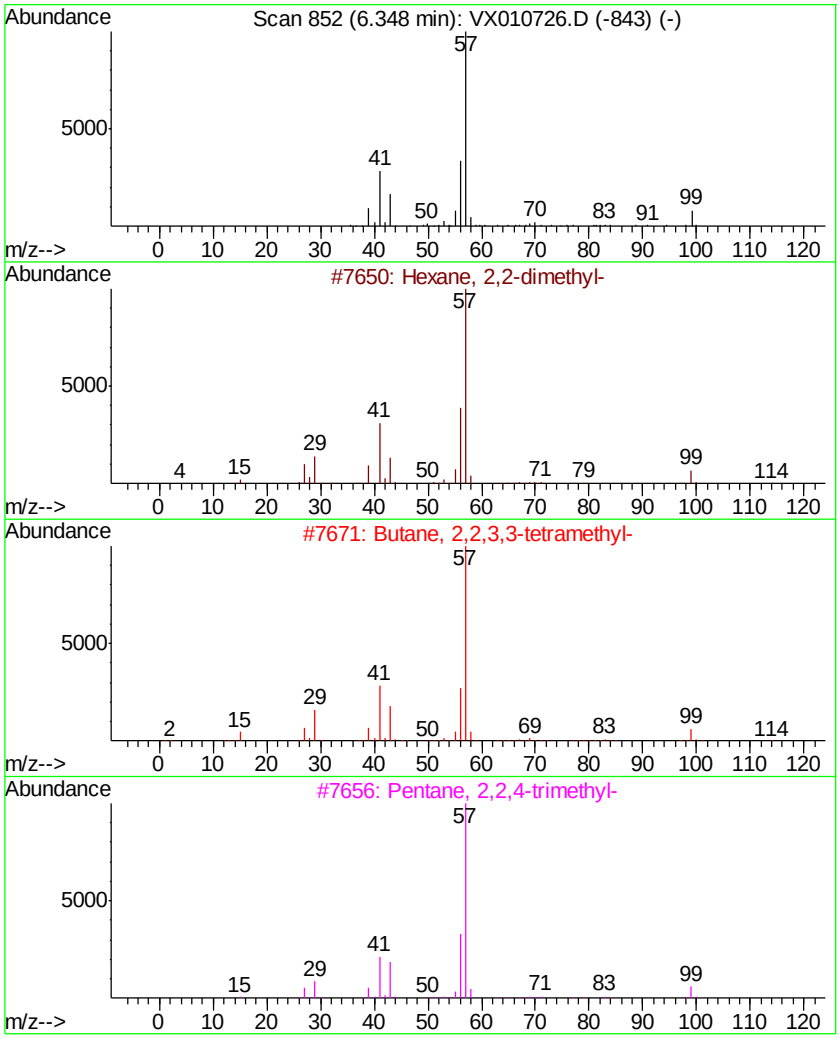
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

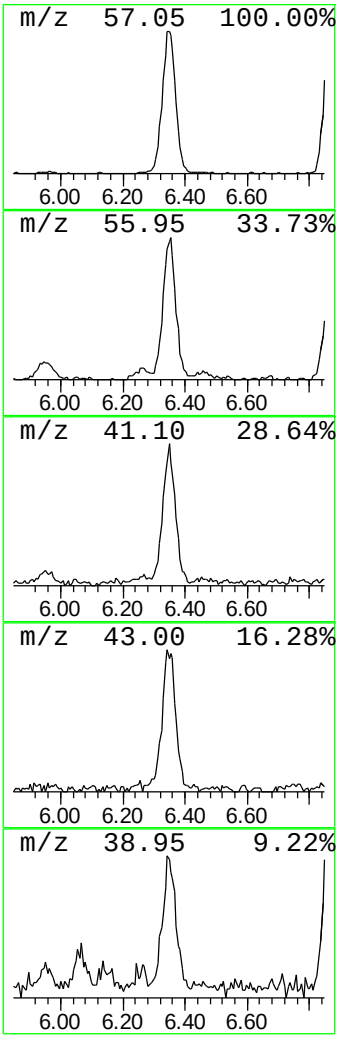
****

Peak Number 3 Hexane, 2,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.35	6.90 ug/l	107623	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
2	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
3	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	64
4	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59





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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

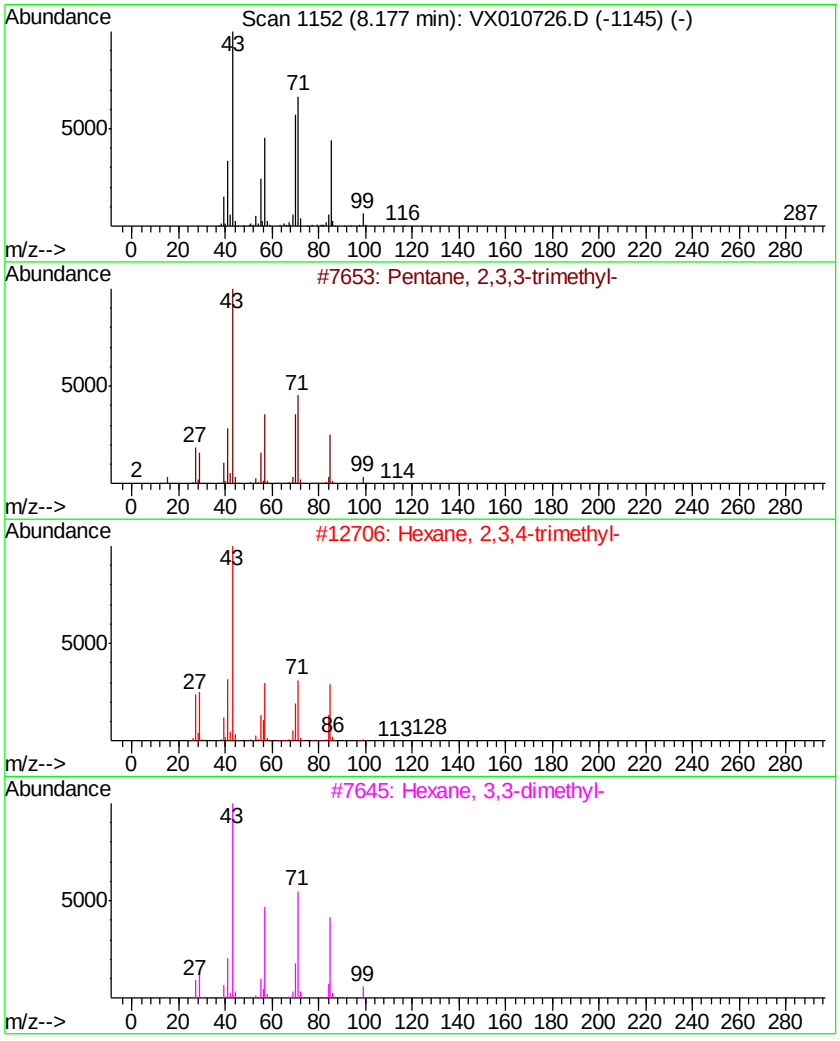
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

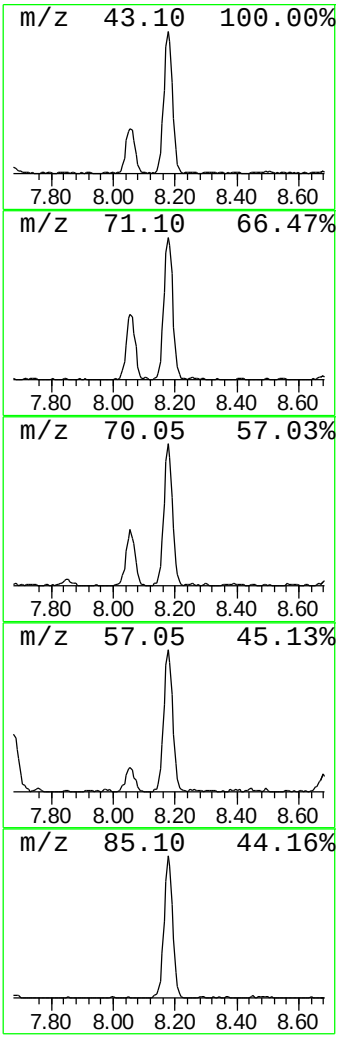
****

Peak Number 4 Pentane, 2,3,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.18	9.90 ug/l	154461	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
3	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
4	1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	53
5	Sulfurous acid, nonyl 2-propyl e...	250	C12H26O3S	1000309-12-0	53





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
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ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_X
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

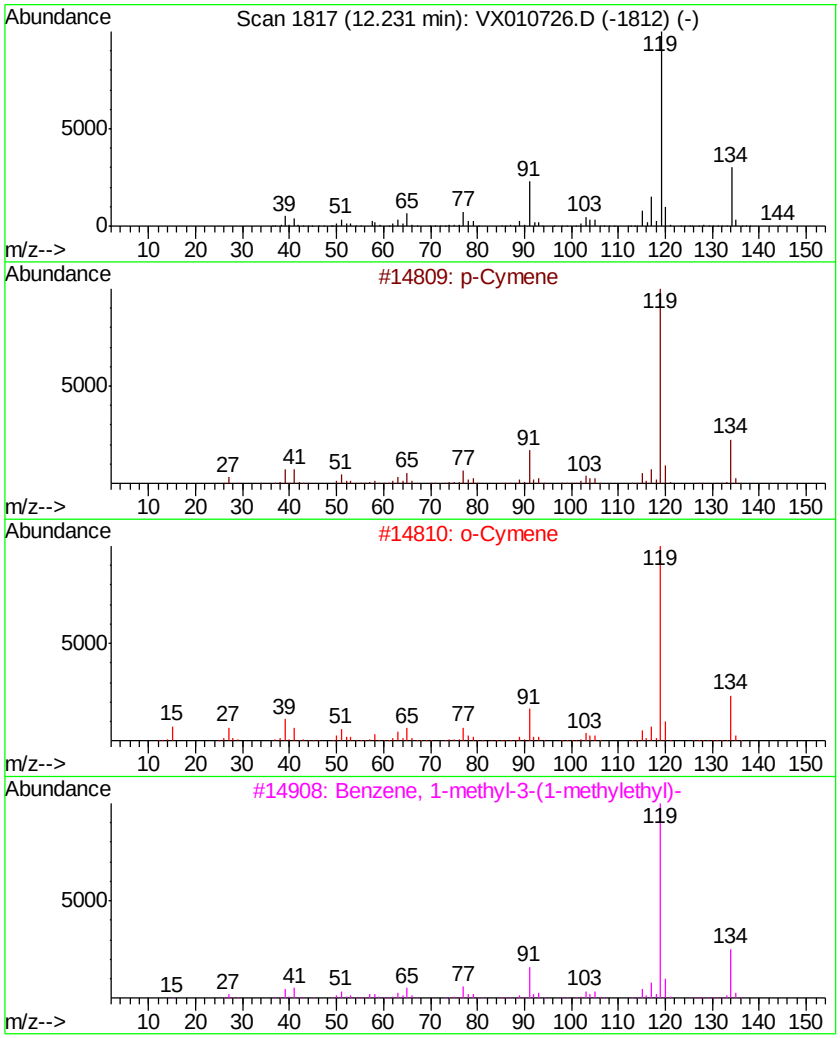
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

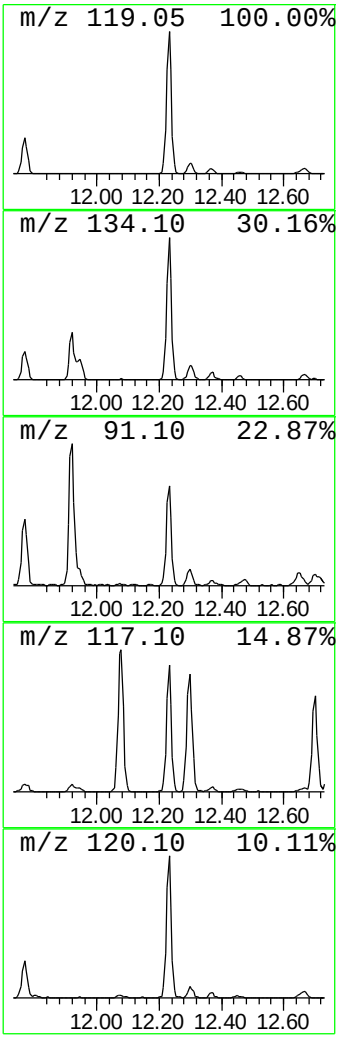
****

Peak Number 5 p-Cymene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.23	21.78 ug/l	410955	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91





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Operator : JC/SP
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ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW082-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

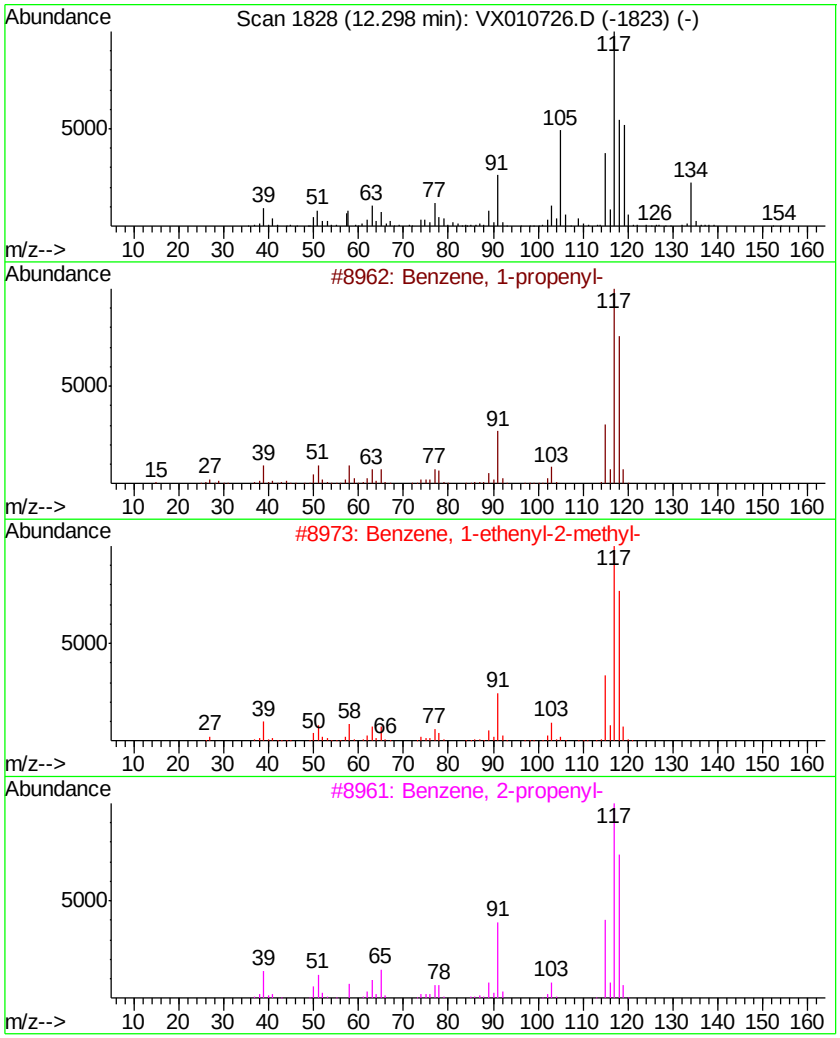
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

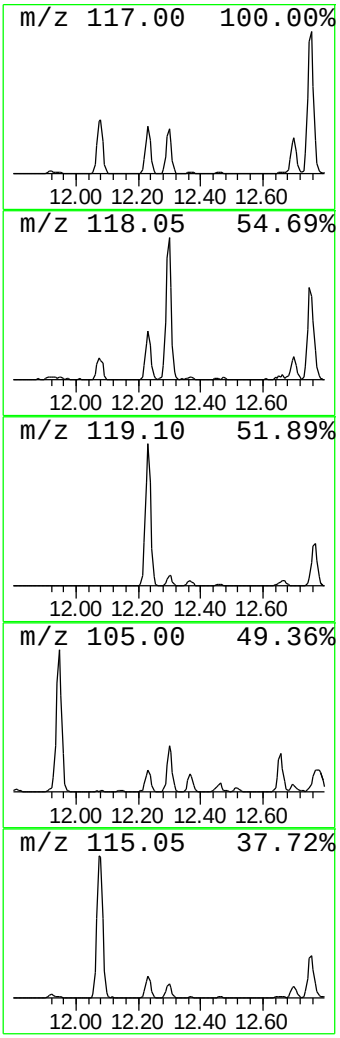
****

Peak Number 6 Benzene, 1-propenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	6.11 ug/l	115232	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propenyl-	118	C9H10	000637-50-3	90
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
4	Indane	118	C9H10	000496-11-7	55
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	55





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010726.D
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Operator : JC/SP
Sample : K3690-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW082-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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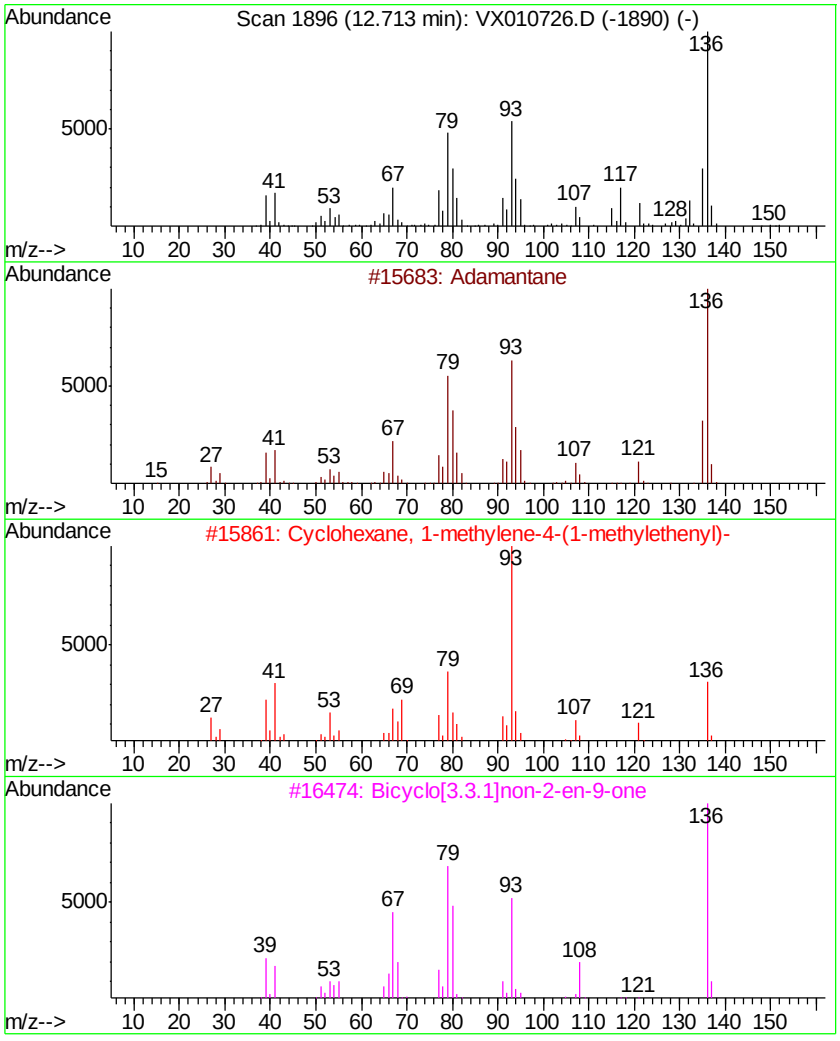
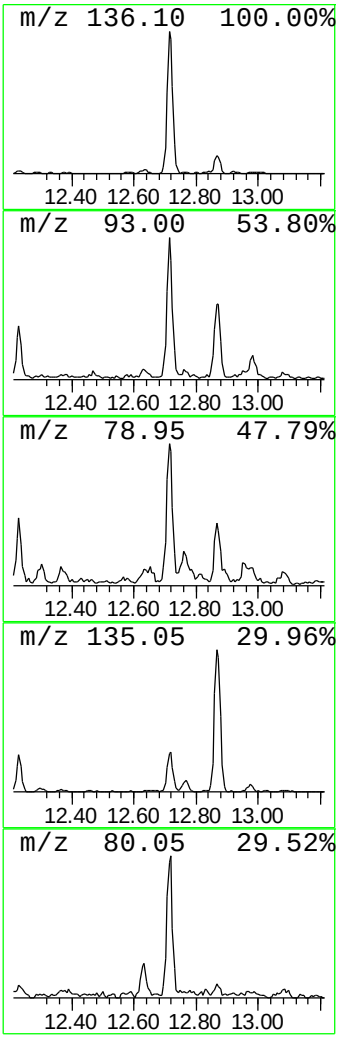
Peak Number 7 Adamantane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.71	8.84 ug/l	166756	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Adamantane	136	C10H16	000281-23-2	98
2	Cyclohexane, 1-methylene-4-(1-me...	136	C10H16	000499-97-8	70
3	Bicyclo[3.3.1]non-2-en-9-one	136	C9H12O	004844-11-5	58
4	.gamma.-Terpinene	136	C10H16	000099-85-4	50
5	1,4-Benzenediamine, 2,3-dimethyl-	136	C8H12N2	005306-96-7	46





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010726.D
Acq On : 09 Jul 2019 03:53
Operator : JC/SP
Sample : K3690-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW082-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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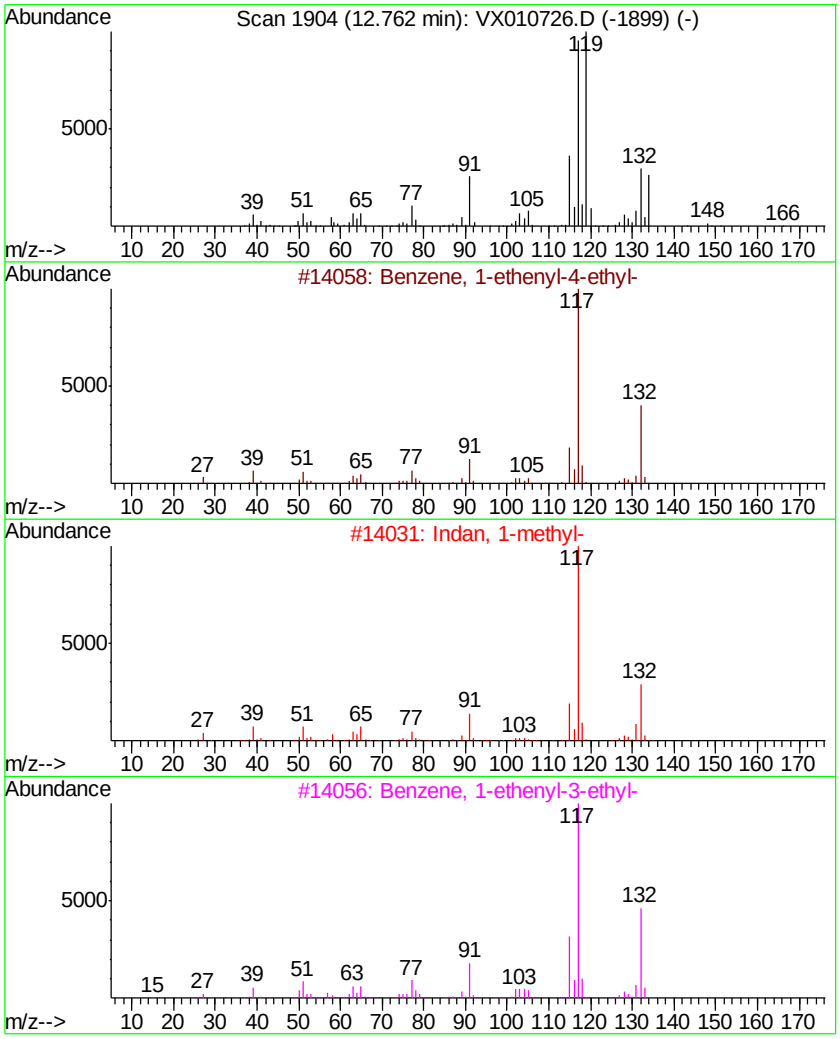
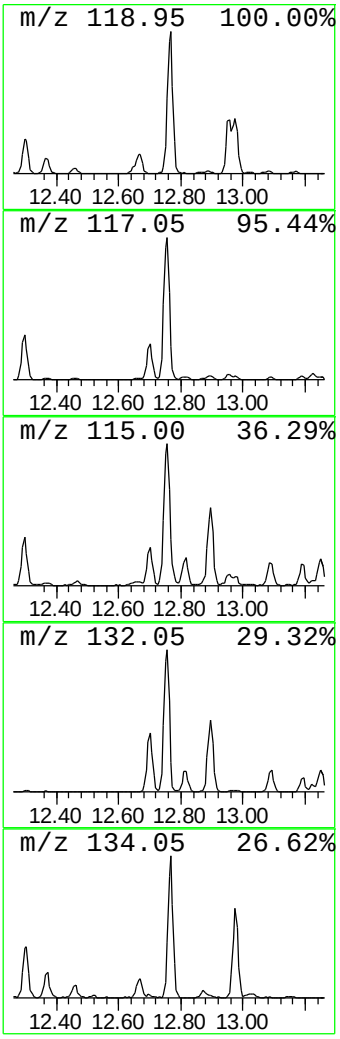
Peak Number 8 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	18.31 ug/l	345427	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
2	Indan, 1-methyl-	132	C10H12	000767-58-8	76
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	74
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	74
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	74





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010726.D
Acq On : 09 Jul 2019 03:53
Operator : JC/SP
Sample : K3690-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW082-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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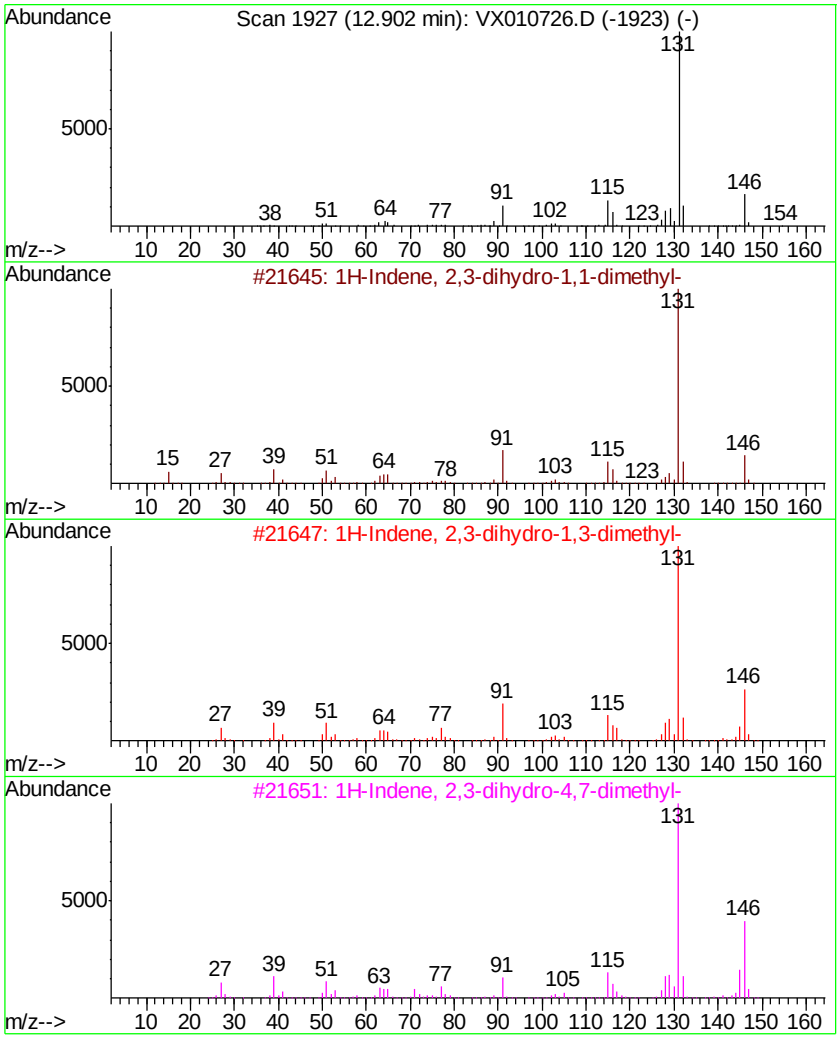
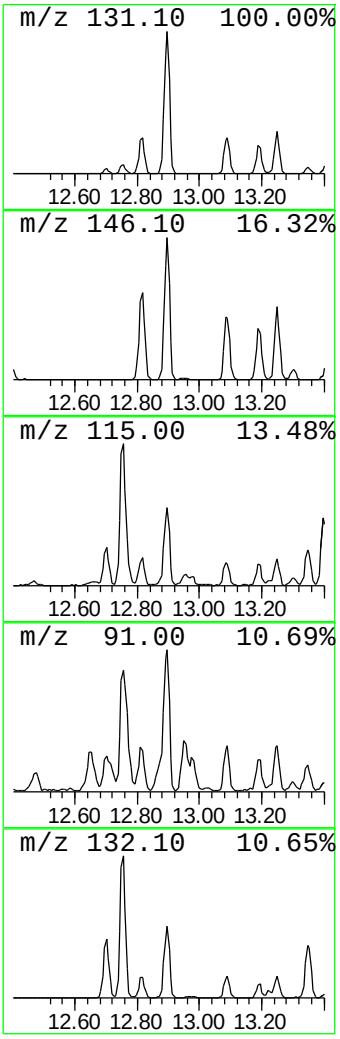
Peak Number 9 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.90	12.29 ug/l	231857	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	94
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	87
5	Cyclopropane, 2-bromo-1-methyl-1...	210	C10H11Br	055091-64-0	78





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010726.D
Acq On : 09 Jul 2019 03:53
Operator : JC/SP
Sample : K3690-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038RW082-190703

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.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
Sulfur dioxide	1.28	5.0	ug/l	65154	1	5.66	650519	50.0
Butane, 2,3-dimet...	2.84	5.8	ug/l	75837	1	5.66	650519	50.0
Hexane, 2,2-dimet...	6.35	6.9	ug/l	107623	2	6.86	779923	50.0
Pentane, 2,3,3-tr...	8.18	9.9	ug/l	154461	2	6.86	779923	50.0
p-Cymene	12.23	21.8	ug/l	410955	4	12.08	943242	50.0
Benzene, 1-propenyl-	12.30	6.1	ug/l	115232	4	12.08	943242	50.0
Adamantane	12.71	8.8	ug/l	166756	4	12.08	943242	50.0
Benzene, 1-etheny...	12.76	18.3	ug/l	345427	4	12.08	943242	50.0
1H-Indene, 2,3-di...	12.90	12.3	ug/l	231857	4	12.08	943242	50.0