LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX080919\
Data File : VX011409.D
Acq On : 09 Aug 2019 14:55
Operator : JC/SP
Sample : K4279-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

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\82X080119W.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	1.782	98	103	113	rVB	134361	188041	1.89%	0.223%
2	1.995	132	138	152	rVB	181173	255503	2.57%	0.304%
3	2.391	195	203	208	rBV	71929	144824	1.46%	0.172%
4	2.836	267	276	279	rBV	141112	306541	3.08%	0.364%
5	2.885	279	284	288	rVV	648845	1372661	13.80%	1.631%
6	2.928	288	291	306	rVB	373340	693110	6.97%	0.824%
7	3.166	319	330	348	rBV	612605	1389324	13.97%	1.651%
8	3.519	378	388	401	rBV	453711	1012802	10.18%	1.204%
9	4.397	522	532	546	rVB	1097759	3179413	31.97%	3.778%
10	5.232	654	669	679	rBV	36837	152429	1.53%	0.181%
11	5.494	700	712	716	rBV	112191	308784	3.11%	0.367%
12	5.580	716	726	735	rBV2	1088100	3477309	34.97%	4.132%
13	5.720	745	749	767	rVB	122655	351094	3.53%	0.417%
14	5.927	770	783	796	rBV4	322487	1012525	10.18%	1.203%
15	6.055	796	804	809	rVV	112976	287623	2.89%	0.342%
16	6.141	809	818	828	rVV	1411974	3700112	37.21%	4.397%
17	6.257	828	837	847	rVV2	150139	521051	5.24%	0.619%
18	6.354	847	853	861	rVV	143098	390804	3.93%	0.464%
19	6.452	861	869	881	rVB2	240969	663859	6.68%	0.789%
20	6.854	926	935	943	rBV	298623	694470	6.98%	0.825%
21	7.464	1021	1035	1046	rBV	1576352	3574455	35.94%	4.248%
22	7.732	1072	1079	1089	rVB	155566	291901	2.94%	0.347%
23	7.848	1089	1098	1106	rVB2	53121	110485	1.11%	0.131%
24	8.024	1121	1127	1139	rVB2	56142	121945	1.23%	0.145%
25	8.342	1174	1179	1182	rBV2	67116	109071	1.10%	0.130%
26	8.384	1182	1186	1192	rVB3	57998	107089	1.08%	0.127%
27	8.500	1200	1205	1214	rBV	81294	150012	1.51%	0.178%
28	8.665	1224	1232	1235	rBV2	216136	390815	3.93%	0.464%
29	8.713	1235	1240	1247	rVB	624132	1055791	10.62%	1.255%
30	8.835	1255	1260	1264	rBV	90273	143717	1.45%	0.171%
31	9.037	1287	1293	1300	rVB	182282	289693	2.91%	0.344%
32	9.165	1304	1314	1319	rBV	86647	145702	1.47%	0.173%
33	9.573	1374	1381	1384	rBV4	82530	140690	1.41%	0.167%
34	9.622	1384	1389	1394	rVB	239689	327805	3.30%	0.390%

LSC Area Percent Report

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Operator : JC/SP
Sample : K4279-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

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\82X080119W.M
Title : SW846 8260

35	9.677	1394	1398	1402	rBV	108774	154923	1.56%	0.184%
36	10.109	1464	1469	1474	rBV	598212	780115	7.84%	0.927%
37	10.183	1474	1481	1486	rVB	87220	132153	1.33%	0.157%
38	10.250	1486	1492	1498	rBV	7695632	9944594	100.00%	11.818%
39	10.359	1501	1510	1516	rVB	2784082	3650178	36.71%	4.338%
40	10.640	1548	1556	1562	rBV3	49948	125644	1.26%	0.149%
41	10.835	1578	1588	1594	rBV2	101374	163452	1.64%	0.194%
42	10.908	1594	1600	1605	rBV4	69812	118087	1.19%	0.140%
43	11.018	1613	1618	1623	rBV	817042	1069510	10.75%	1.271%
44	11.134	1631	1637	1642	rBV	555843	693512	6.97%	0.824%
45	11.359	1669	1674	1680	rBV	984930	1216403	12.23%	1.446%
46	11.451	1682	1689	1693	rVV	1383305	1813454	18.24%	2.155%
47	11.506	1693	1698	1707	rVB	1054410	1296093	13.03%	1.540%
48	11.664	1714	1724	1734	rBV	3659323	4643212	46.69%	5.518%
49	11.804	1742	1747	1752	rBV	7877343	9533814	95.87%	11.330%
50	11.945	1767	1770	1775	rVV	191749	234416	2.36%	0.279%
51	12.024	1776	1783	1786	rVV	146386	184286	1.85%	0.219%
52	12.073	1786	1791	1797	rVB2	775770	1128288	11.35%	1.341%
53	12.140	1797	1802	1807	rBV	3155299	3741809	37.63%	4.447%
54	12.231	1812	1817	1821	rBV	89210	109931	1.11%	0.131%
55	12.298	1823	1828	1835	rBV	1192973	1628805	16.38%	1.936%
56	12.365	1835	1839	1849	rVB3	604229	986696	9.92%	1.173%
57	12.457	1849	1854	1859	rBV2	246461	317720	3.19%	0.378%
58	12.518	1859	1864	1869	rVB	642754	780908	7.85%	0.928%
59	12.585	1870	1875	1877	rBV	615362	808674	8.13%	0.961%
60	12.609	1877	1879	1883	rVB	713275	706386	7.10%	0.839%
61	12.664	1883	1888	1892	rBV	745196	875515	8.80%	1.040%
62	12.700	1892	1894	1898	rVB	164877	165686	1.67%	0.197%
63	12.755	1898	1903	1910	rBV4	468052	864667	8.69%	1.028%
64	12.902	1922	1927	1931	rVB2	469811	605777	6.09%	0.720%
65	12.975	1931	1939	1942	rBV	429584	561168	5.64%	0.667%
66	13.018	1942	1946	1952	rVB	773488	940297	9.46%	1.117%
67	13.078	1952	1956	1961	rBV3	100795	177645	1.79%	0.211%
68	13.194	1971	1975	1977	rBV2	116054	145741	1.47%	0.173%
69	13.219	1977	1979	1983	rVB	141111	165697	1.67%	0.197%
70	13.304	1990	1993	1995	rBV	118989	145613	1.46%	0.173%
71	13.341	1995	1999	2005	rVB2	1240840	1684400	16.94%	2.002%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX080919\
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Acq On : 09 Aug 2019 14:55
Operator : JC/SP
Sample : K4279-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

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\82X080119W.M
Title : SW846 8260

72	13.475	2017	2021	2029	rVB	665215	871832	8.77%	1.036%
73	13.560	2029	2035	2038	rBV3	67306	104837	1.05%	0.125%
74	13.639	2044	2048	2053	rVV3	148984	271263	2.73%	0.322%
75	13.719	2059	2061	2067	rVB2	157263	214170	2.15%	0.255%
76	13.834	2072	2080	2087	rVB	1025956	1420336	14.28%	1.688%
77	13.908	2087	2092	2097	rVB	112706	154327	1.55%	0.183%
78	13.981	2097	2104	2108	rBV2	155327	239204	2.41%	0.284%
79	14.286	2147	2154	2159	rVB2	206238	363659	3.66%	0.432%
80	14.536	2190	2195	2200	rBV2	210591	286599	2.88%	0.341%
81	14.694	2214	2221	2225	rBV	330718	444868	4.47%	0.529%
82	14.834	2239	2244	2249	rBV	325527	421921	4.24%	0.501%

Sum of corrected areas: 84149735

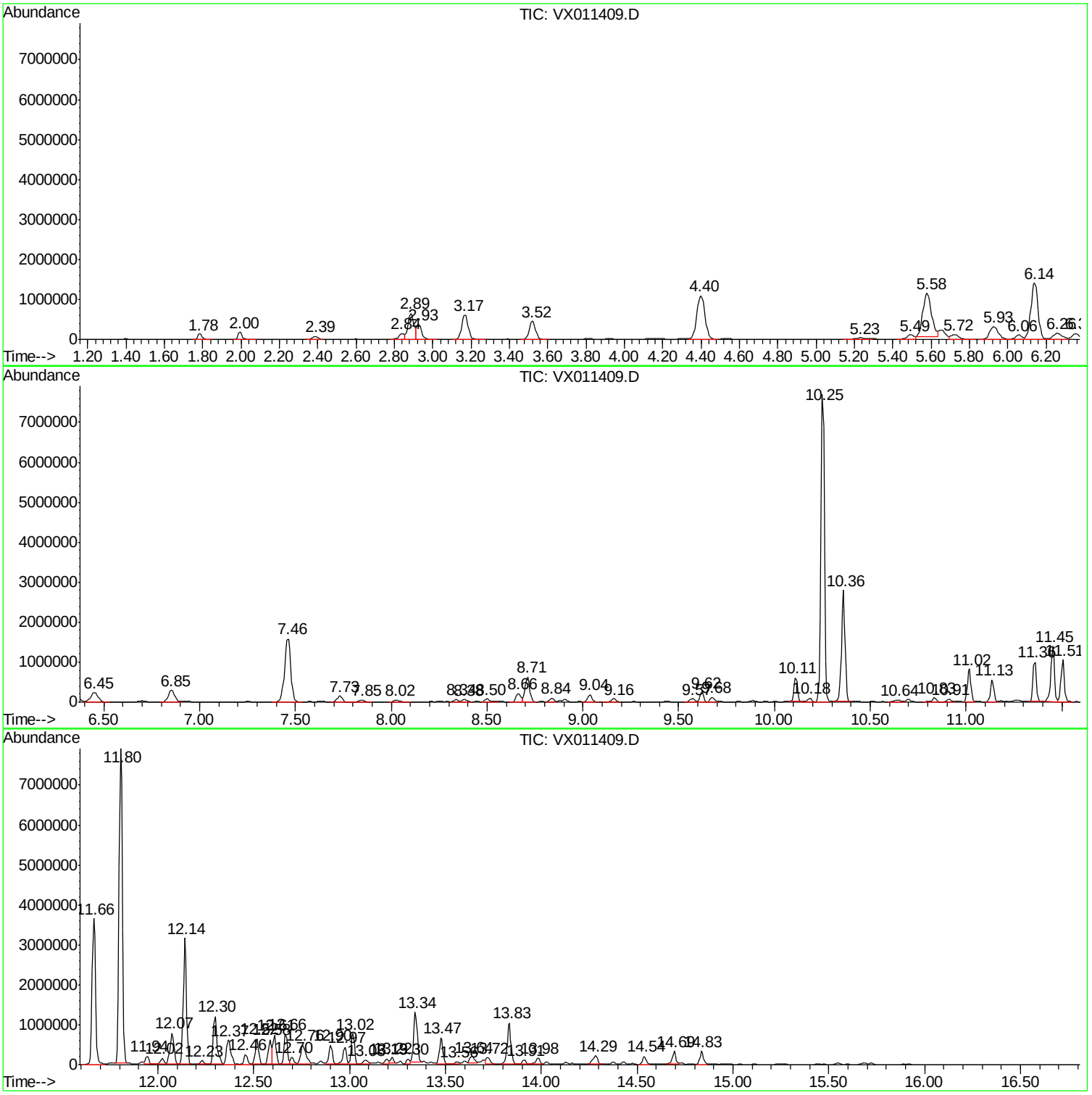
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX080919\
Data File : VX011409.D
Acq On : 09 Aug 2019 14:55
Operator : JC/SP
Sample : K4279-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.M
Quant Title : SW846 8260

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



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX080919\
Data File : VX011409.D
Acq On : 09 Aug 2019 14:55
Operator : JC/SP
Sample : K4279-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.M
Quant Title : SW846 8260

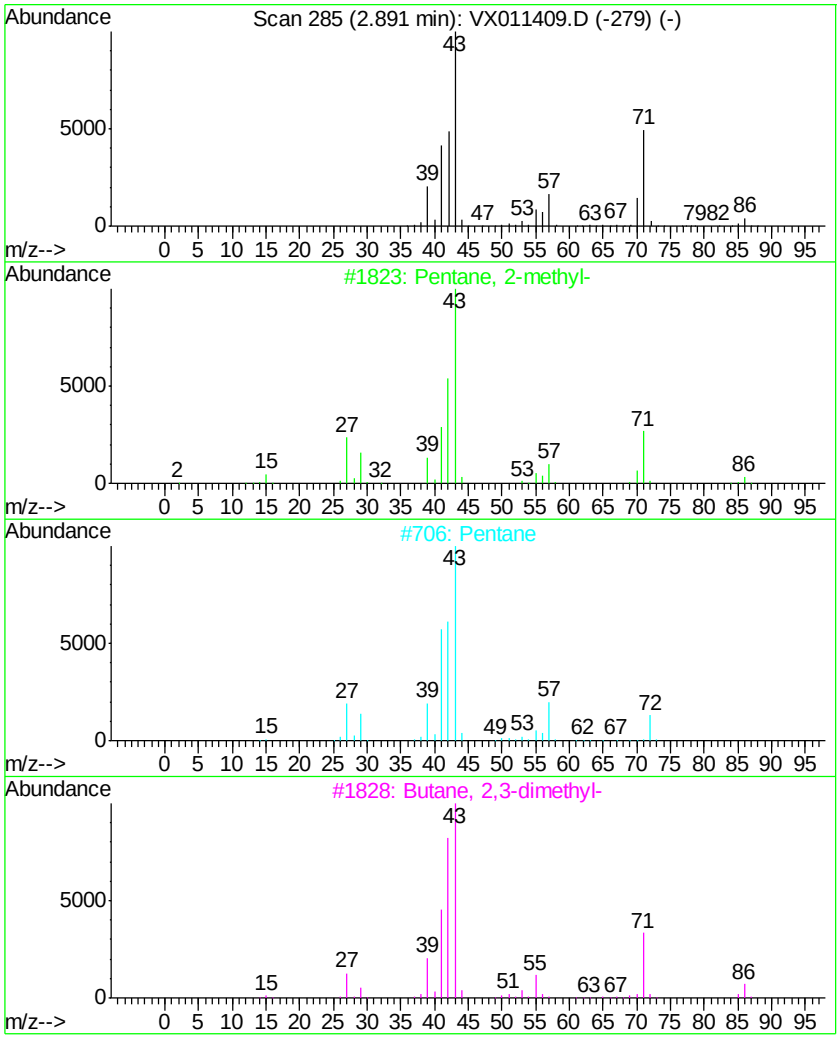
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

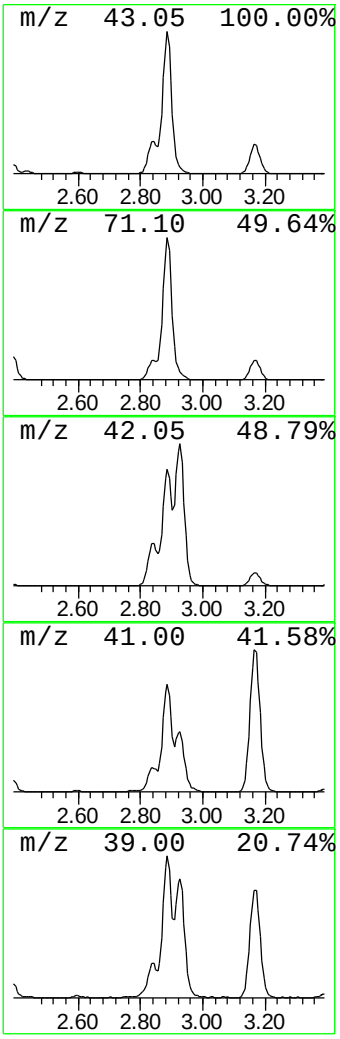
****

Peak Number 1 Pentane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.89	195.48 ug/l	1372660	Pentafluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2	Pentane	72	C5H12	000109-66-0	46
3	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	46
4	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45
5	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX080919\
Data File : VX011409.D
Acq On : 09 Aug 2019 14:55
Operator : JC/SP
Sample : K4279-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.M
Quant Title : SW846 8260

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

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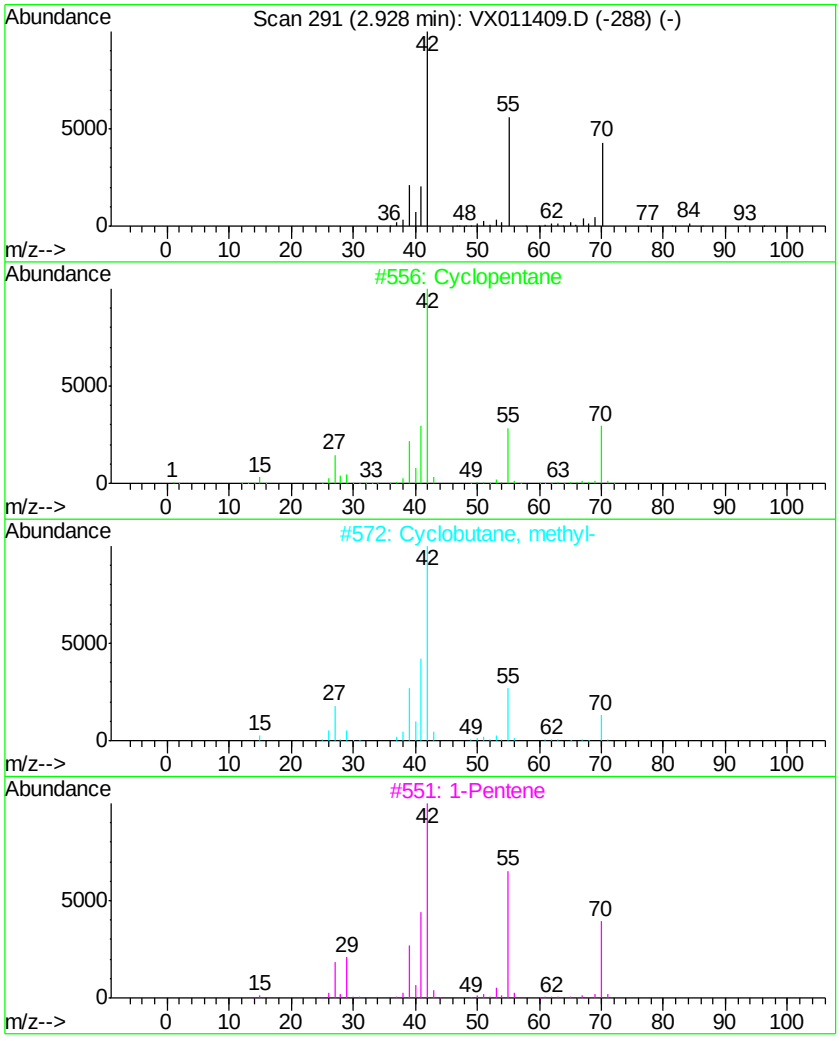
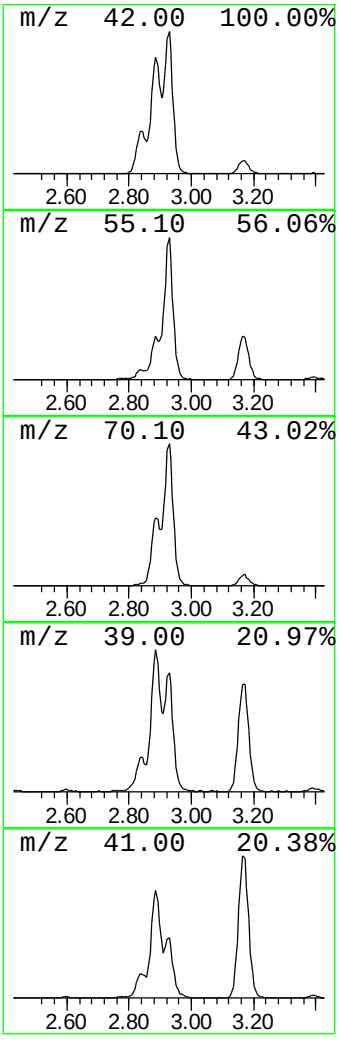
Peak Number 2 Cyclopentane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.93	98.71 ug/l	693110	Pentafluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane	70	C5H10	000287-92-3	78
2	Cyclobutane, methyl-	70	C5H10	000598-61-8	72
3	1-Pentene	70	C5H10	000109-67-1	56
4	2-Pentene, (Z)-	70	C5H10	000627-20-3	35
5	2-Pentene	70	C5H10	000109-68-2	25





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX080919\
Data File : VX011409.D
Acq On : 09 Aug 2019 14:55
Operator : JC/SP
Sample : K4279-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.M
Quant Title : SW846 8260

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

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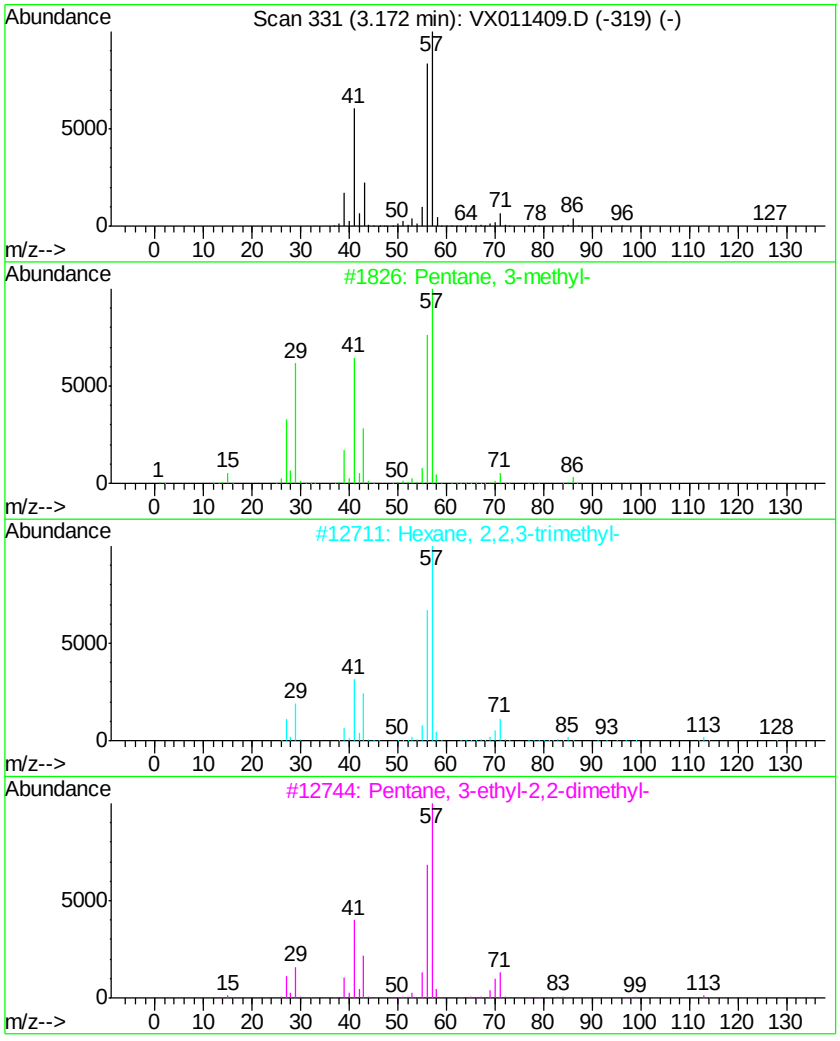
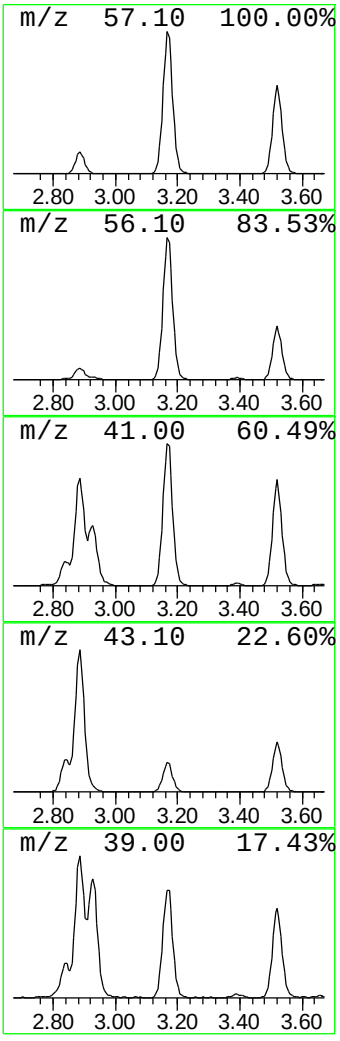
Peak Number 3 Pentane, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.17	197.86 ug/l	1389320	Pentafluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	83
3	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
4	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45
5	Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	40





Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX080919\
Data File : VX011409.D
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Operator : JC/SP
Sample : K4279-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.M
Quant Title : SW846 8260

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

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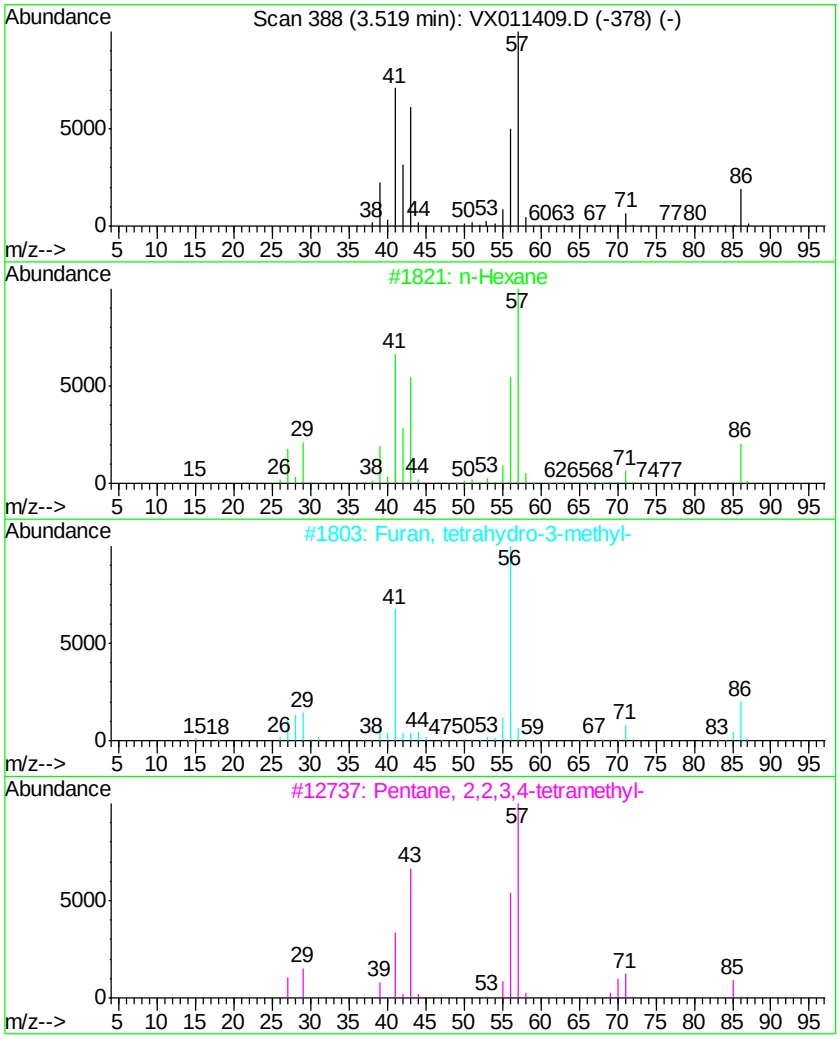
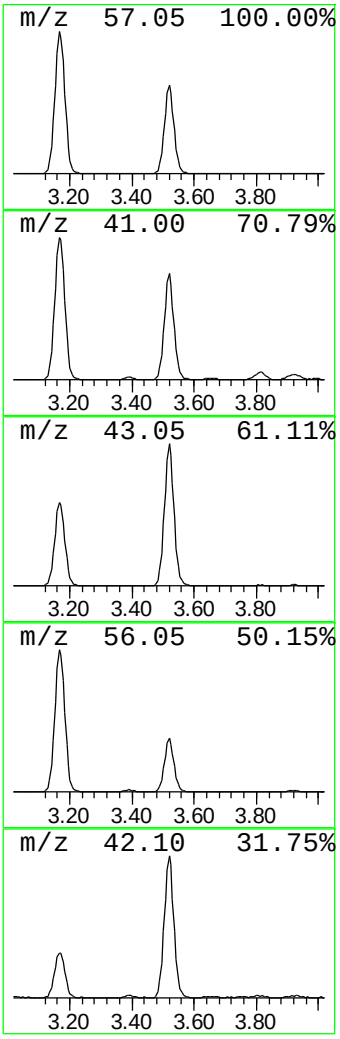
Peak Number 4 n-Hexane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.52	144.24 ug/l	1012800	Pentafluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	94
2	Furan, tetrahydro-3-methyl-	86	C5H10O	013423-15-9	53
3	Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	50
4	Hydroperoxide, hexyl	118	C6H14O2	004312-76-9	38
5	Butanal, 2-methyl-	86	C5H10O	000096-17-3	38





Library Search Compound Report

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Sample : K4279-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.M
Quant Title : SW846 8260

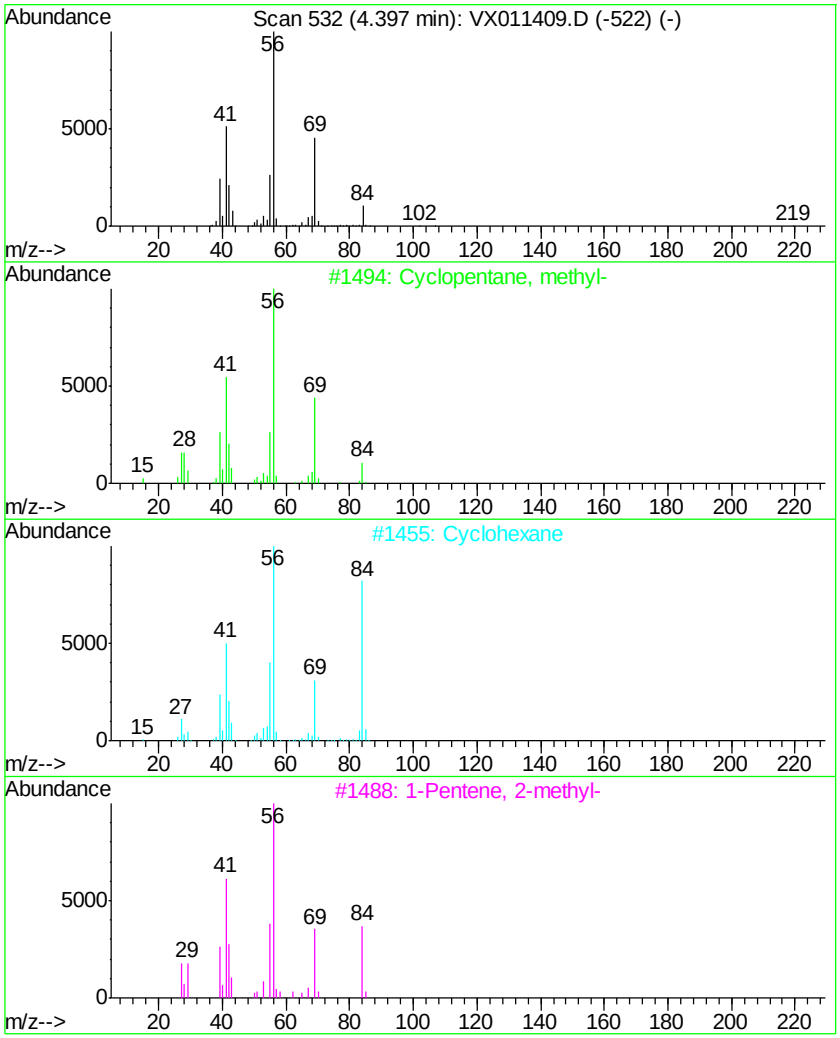
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

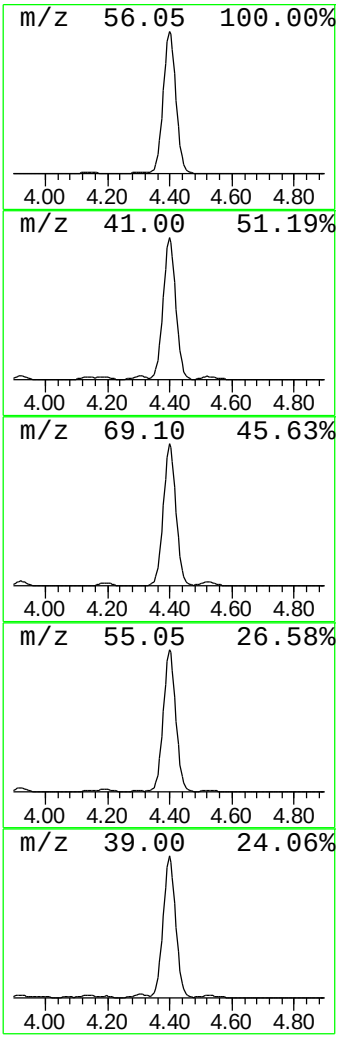
****

Peak Number 5 Cyclopentane, methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	452.79 ug/l	3179410	Pentafluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Cyclohexane	84	C6H12	000110-82-7	80
3	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72
4	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX080919\
Data File : VX011409.D
Acq On : 09 Aug 2019 14:55
Operator : JC/SP
Sample : K4279-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.M
Quant Title : SW846 8260

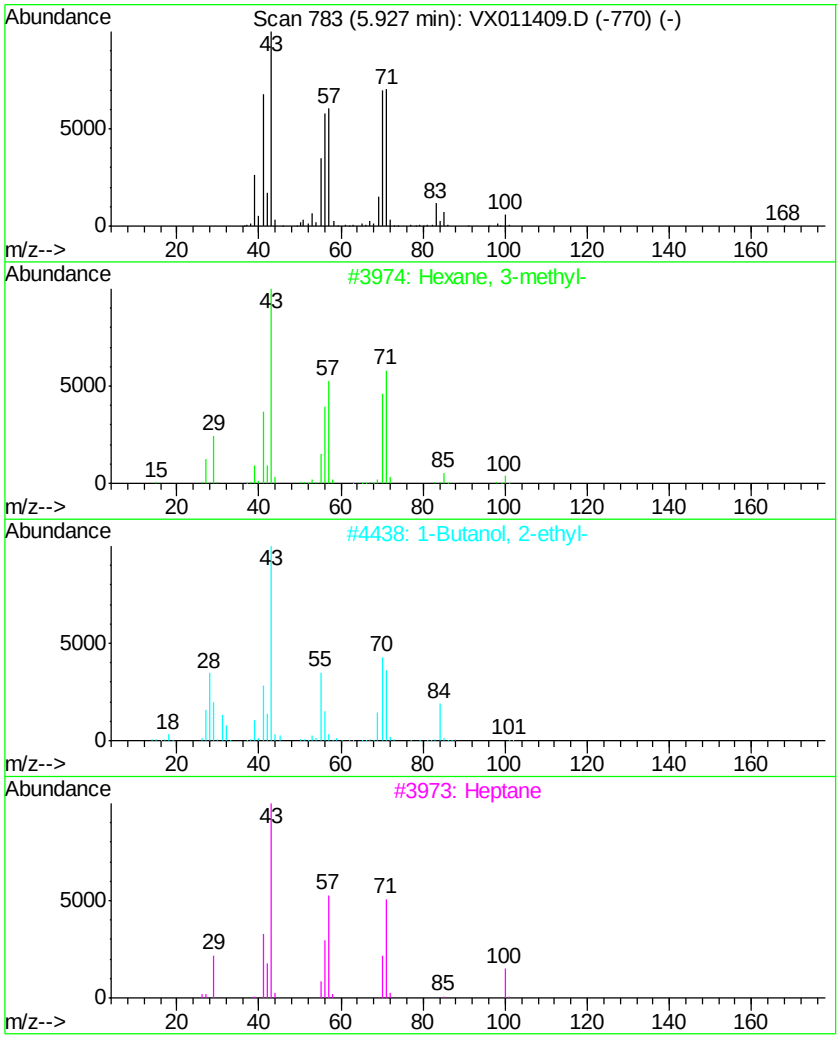
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

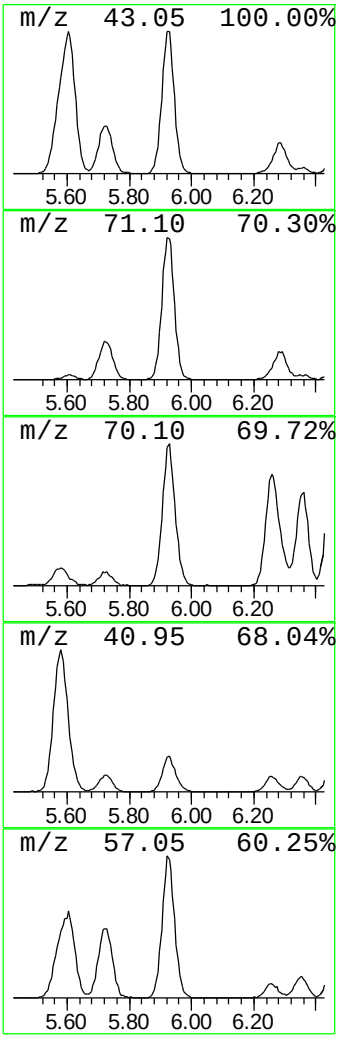
****

Peak Number 6 Hexane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.93	144.20 ug/l	1012530	Pentafluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-	100	C7H16	000589-34-4	86
2	1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	53
3	Heptane	100	C7H16	000142-82-5	52
4	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	50
5	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	50





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX080919\
Data File : VX011409.D
Acq On : 09 Aug 2019 14:55
Operator : JC/SP
Sample : K4279-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.M
Quant Title : SW846 8260

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

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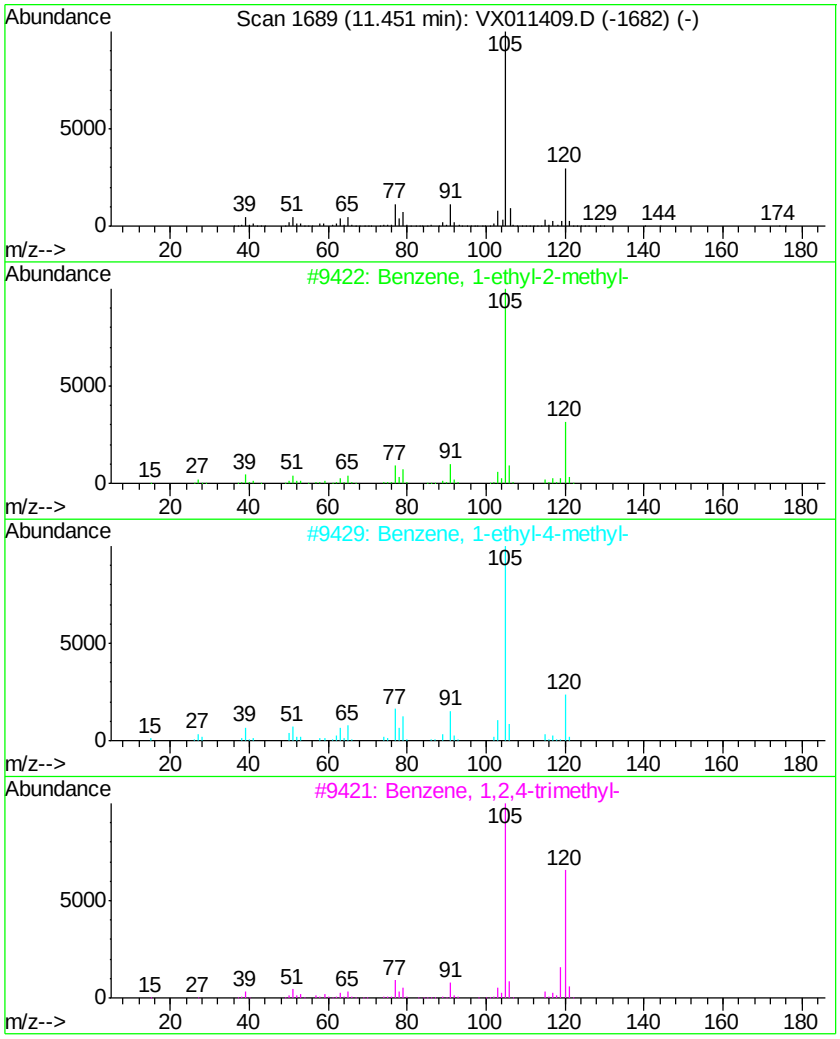
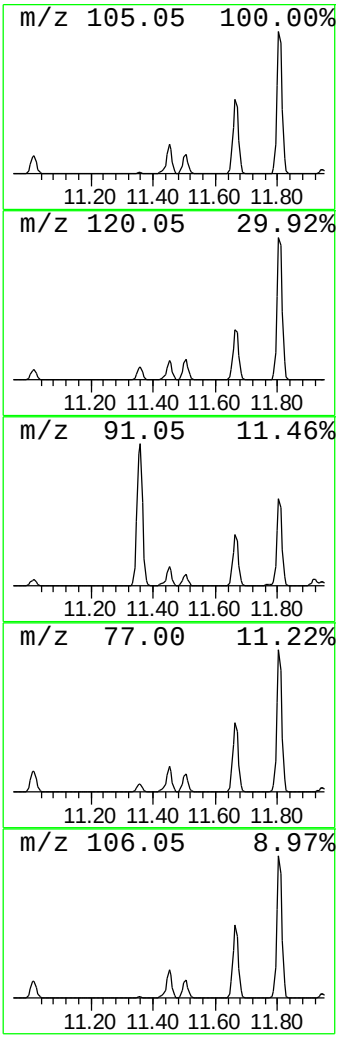
Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.45	80.36 ug/l	1813450	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	95
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX080919\
Data File : VX011409.D
Acq On : 09 Aug 2019 14:55
Operator : JC/SP
Sample : K4279-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.M
Quant Title : SW846 8260

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

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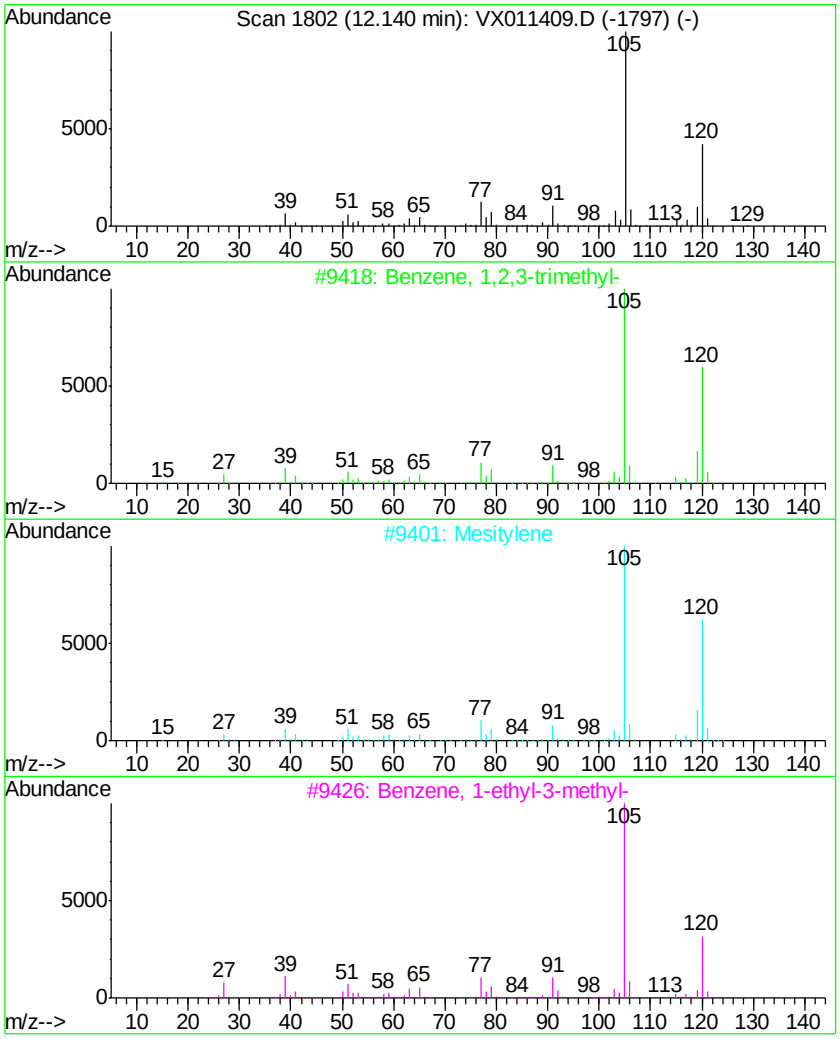
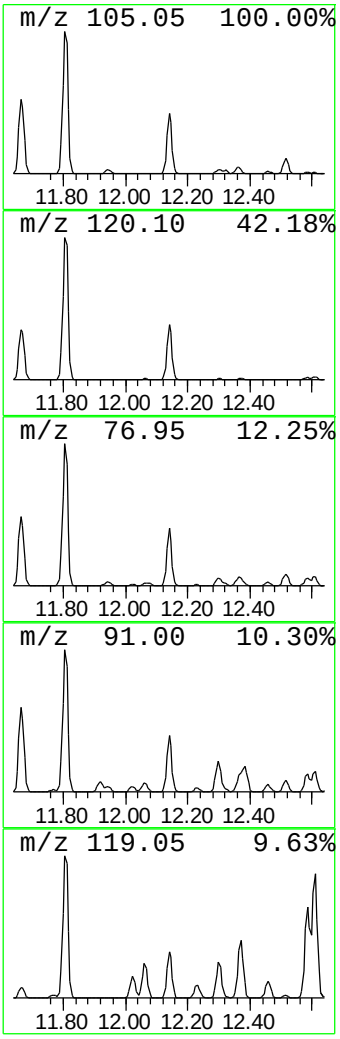
Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	165.82 ug/l	3741810	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	91
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX080919\
Data File : VX011409.D
Acq On : 09 Aug 2019 14:55
Operator : JC/SP
Sample : K4279-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.M
Quant Title : SW846 8260

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

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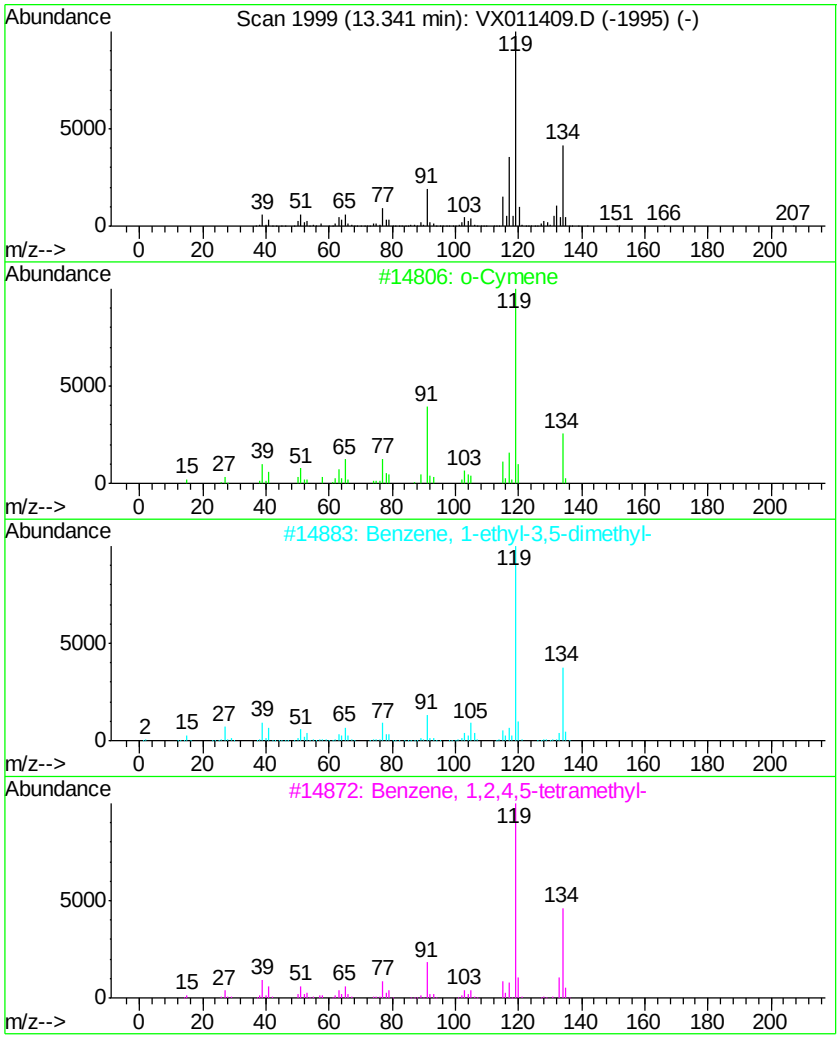
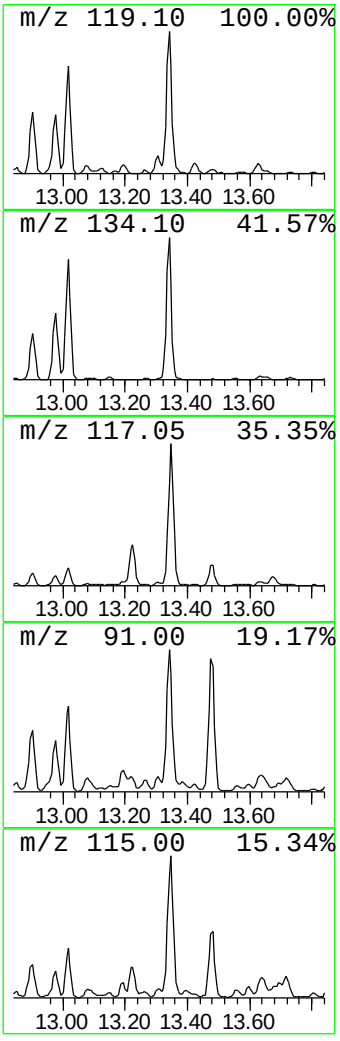
Peak Number 10 o-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	74.64 ug/l	1684400	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	81
2	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	81
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	76
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	76
5	p-Cymene	134	C10H14	000099-87-6	74





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX080919\
Data File : VX011409.D
Acq On : 09 Aug 2019 14:55
Operator : JC/SP
Sample : K4279-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : Z3-0671

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.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
Pentane, 2-methyl-	2.89	195.5	ug/l	1372660	1	5.66	351094	50.0
Cyclopentane	2.93	98.7	ug/l	693110	1	5.66	351094	50.0
Pentane, 3-methyl-	3.17	197.9	ug/l	1389320	1	5.66	351094	50.0
n-Hexane	3.52	144.2	ug/l	1012800	1	5.66	351094	50.0
Cyclopentane, met...	4.40	452.8	ug/l	3179410	1	5.66	351094	50.0
Hexane, 3-methyl-	5.93	144.2	ug/l	1012530	1	5.66	351094	50.0
Benzene, 1-ethyl-...	11.45	80.4	ug/l	1813450	4	12.08	1128290	50.0
Benzene, 1,2,3-tr...	12.14	165.8	ug/l	3741810	4	12.08	1128290	50.0
o-Cymene	13.34	74.6	ug/l	1684400	4	12.08	1128290	50.0