LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW102819\
Data File : VW013780.D
Acq On : 28 Oct 2019 21:09
Operator : SY/VA
Sample : K5563-11
Misc : 7.10G/5ML/MSVOA W/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : TP-2

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_W\METHOD\82W102319S.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	7.885	971	981	986	rBV	153995	371740	31.12%	2.468%
2	7.946	986	991	1001	rVB	291512	641666	53.72%	4.260%
3	8.305	1040	1050	1062	rBV	136780	304339	25.48%	2.021%
4	8.842	1129	1138	1151	rBV	408966	812317	68.01%	5.394%
5	10.323	1373	1381	1390	rBV	619078	1092996	91.51%	7.257%
6	10.488	1403	1408	1416	rVB5	7253	17329	1.45%	0.115%
7	10.665	1431	1437	1441	rBV2	20084	34224	2.87%	0.227%
8	10.756	1447	1452	1458	rBV5	8985	17795	1.49%	0.118%
9	10.933	1477	1481	1488	rVB2	9211	16729	1.40%	0.111%
10	11.036	1491	1498	1500	rBV5	6652	13546	1.13%	0.090%
11	11.110	1505	1510	1513	rBV3	12526	23803	1.99%	0.158%
12	11.140	1513	1515	1518	rVV3	15658	23944	2.00%	0.159%
13	11.177	1518	1521	1524	rVV2	16438	25533	2.14%	0.170%
14	11.232	1524	1530	1535	rVV	158379	279204	23.38%	1.854%
15	11.286	1535	1539	1542	rVV2	22678	37815	3.17%	0.251%
16	11.335	1542	1547	1551	rVB4	24966	44338	3.71%	0.294%
17	11.433	1557	1563	1571	rVB	78476	143201	11.99%	0.951%
18	11.512	1571	1576	1580	rBV2	12984	22790	1.91%	0.151%
19	11.628	1588	1595	1605	rBV	493159	855657	71.64%	5.681%
20	11.725	1606	1611	1613	rBV3	18183	29632	2.48%	0.197%
21	11.762	1613	1617	1621	rVV2	68933	94110	7.88%	0.625%
22	11.817	1621	1626	1631	rBV2	69216	130797	10.95%	0.868%
23	11.896	1631	1639	1647	rVB4	134171	458669	38.40%	3.045%
24	11.969	1647	1651	1657	rBV4	34625	71751	6.01%	0.476%
25	12.036	1657	1662	1670	rVB3	64011	127912	10.71%	0.849%
26	12.134	1670	1678	1685	rBV2	212751	542547	45.42%	3.602%
27	12.195	1685	1688	1690	rBV2	31115	41197	3.45%	0.274%
28	12.225	1690	1693	1695	rVV2	36953	54302	4.55%	0.361%
29	12.250	1695	1697	1701	rVB	86200	103427	8.66%	0.687%
30	12.298	1701	1705	1706	rBV2	57569	72237	6.05%	0.480%
31	12.341	1706	1712	1722	rVB5	214595	537740	45.02%	3.570%
32	12.451	1722	1730	1737	rBV9	90389	287368	24.06%	1.908%
33	12.561	1746	1748	1752	rVB3	58397	67412	5.64%	0.448%
34	12.615	1752	1757	1763	rBV	411396	677042	56.68%	4.495%

LSC Area Percent Report

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Integration Parameters: RTEINT.P
Integrator: RTE
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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_W\METHOD\82W102319S.M
Title : SW846 8260

35	12.695	1763	1770	1775	rBV4	135346	334620	28.01%	2.222%
36	12.780	1778	1784	1791	rVB3	640647	1194448	100.00%	7.931%
37	12.859	1791	1797	1803	rBV6	154864	431987	36.17%	2.868%
38	12.945	1804	1811	1823	rVB8	238953	829910	69.48%	5.510%
39	13.079	1829	1833	1837	rVB3	156388	226410	18.96%	1.503%
40	13.121	1837	1840	1843	rVB3	25156	27297	2.29%	0.181%
41	13.170	1843	1848	1857	rBV7	135452	403856	33.81%	2.681%
42	13.256	1857	1862	1866	rBV2	137629	246491	20.64%	1.637%
43	13.378	1878	1882	1886	rBV3	85270	139549	11.68%	0.927%
44	13.432	1887	1891	1896	rVV5	92737	193462	16.20%	1.285%
45	13.481	1896	1899	1904	rVB5	49111	67521	5.65%	0.448%
46	13.554	1906	1911	1919	rVB	381700	704390	58.97%	4.677%
47	13.652	1919	1927	1930	rBV2	62272	161200	13.50%	1.070%
48	13.701	1930	1935	1942	rVB4	94822	188027	15.74%	1.248%
49	13.816	1948	1954	1958	rBV4	34991	82699	6.92%	0.549%
50	13.877	1958	1964	1970	rVB3	234054	413898	34.65%	2.748%
51	14.012	1978	1986	1989	rBV6	74743	138085	11.56%	0.917%
52	14.097	1995	2000	2006	rVB	147559	256554	21.48%	1.703%
53	14.158	2006	2010	2012	rVV2	33708	54017	4.52%	0.359%
54	14.207	2012	2018	2023	rVV2	180031	340156	28.48%	2.259%
55	14.268	2023	2028	2034	rVB6	41245	82554	6.91%	0.548%
56	14.341	2034	2040	2044	rBV	45223	77074	6.45%	0.512%
57	14.383	2044	2047	2050	rVV4	28091	46363	3.88%	0.308%
58	14.420	2050	2053	2056	rVV2	46836	72920	6.10%	0.484%
59	14.457	2056	2059	2063	rVB	42331	57695	4.83%	0.383%
60	14.536	2068	2072	2081	rVB7	17305	51838	4.34%	0.344%
61	14.639	2086	2089	2094	rVB2	21900	33753	2.83%	0.224%
62	14.700	2094	2099	2101	rBV3	12464	21020	1.76%	0.140%
63	14.719	2101	2102	2109	rVB5	15782	22333	1.87%	0.148%
64	14.798	2109	2115	2122	rVB2	38343	85768	7.18%	0.569%

Sum of corrected areas: 15061004

LSC Report - Integrated Chromatogram

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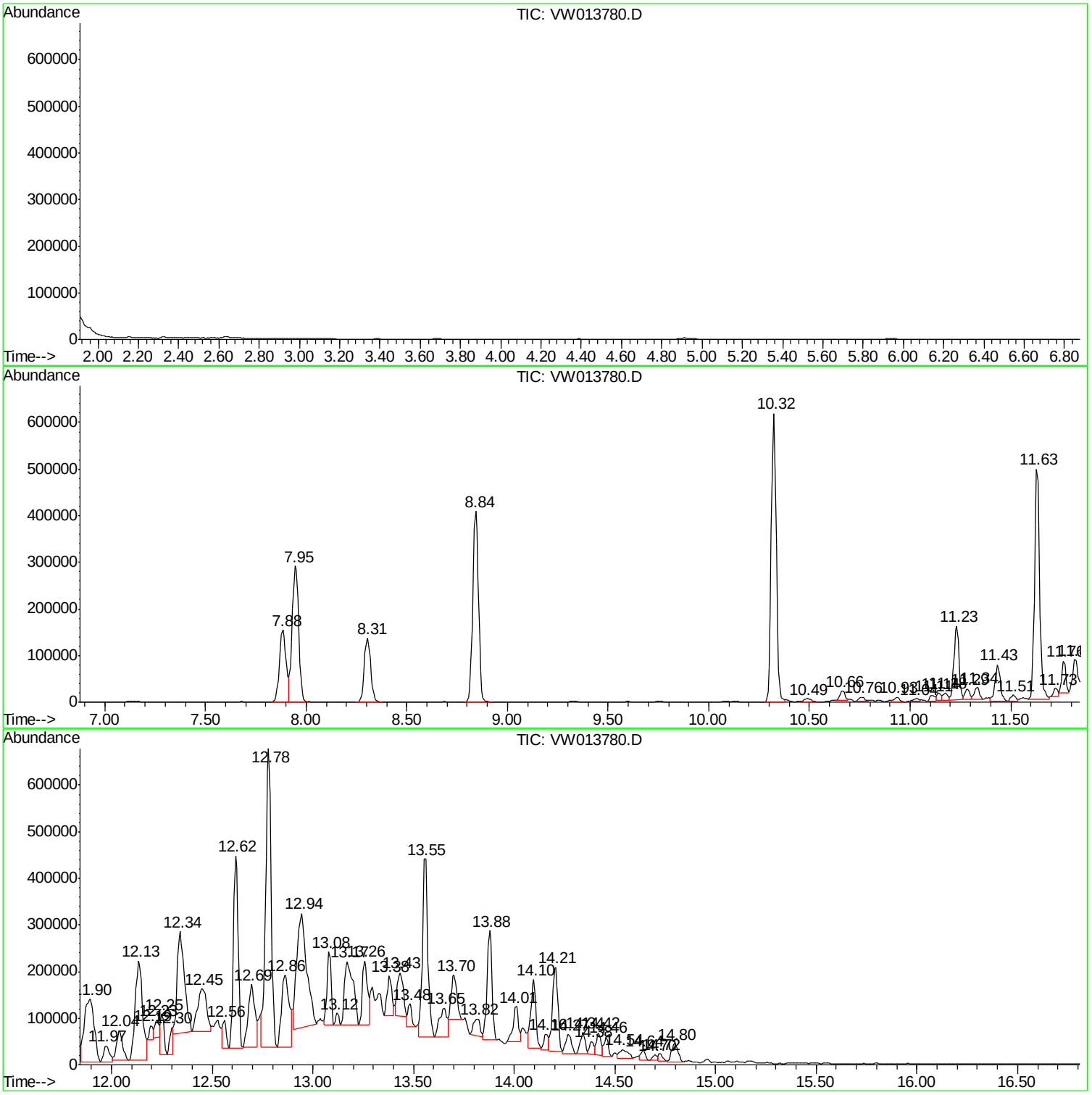
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : TP-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W102319S.M
Quant Title : SW846 8260

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



Library Search Compound Report

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ALS Vial : 23 Sample Multiplier: 1

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Quant Title : SW846 8260

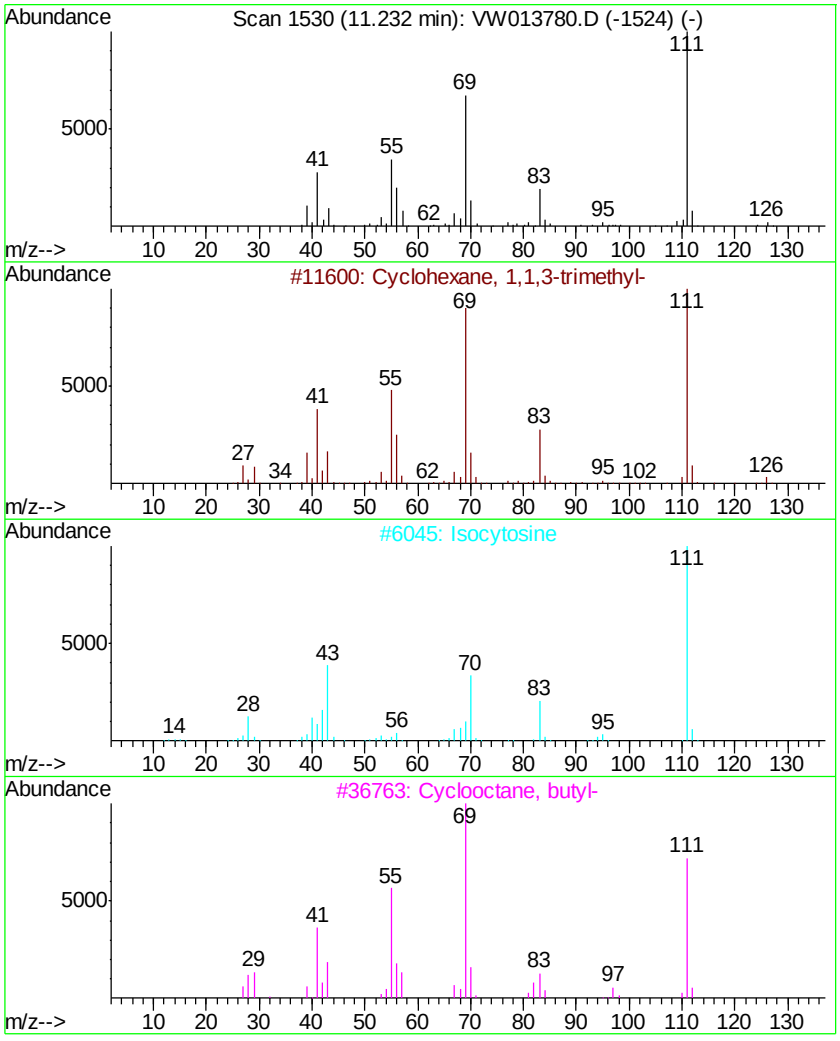
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

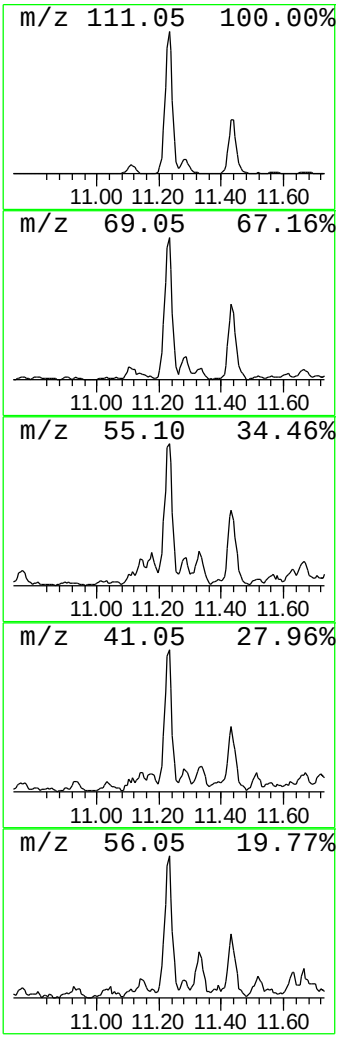
****

Peak Number 1 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.23	16.32 ug/l	279204	Chlorobenzene-d5	11.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	97
2	Isocytosine	111	C4H5N3O	000108-53-2	59
3	Cyclooctane, butyl-	168	C12H24	016538-93-5	56
4	Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	53
5	Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	53





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Data File : VW013780.D
Acq On : 28 Oct 2019 21:09
Operator : SY/VA
Sample : K5563-11
Misc : 7.10G/5ML/MSVOA W/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : TP-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W102319S.M
Quant Title : SW846 8260

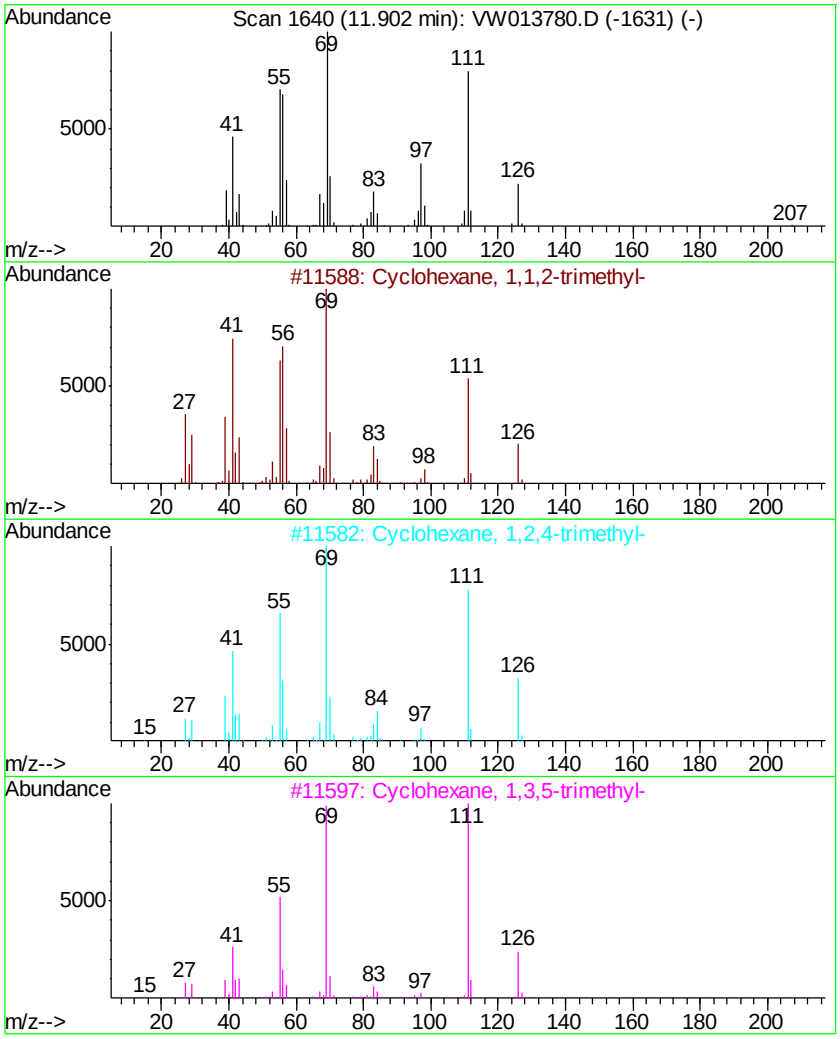
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

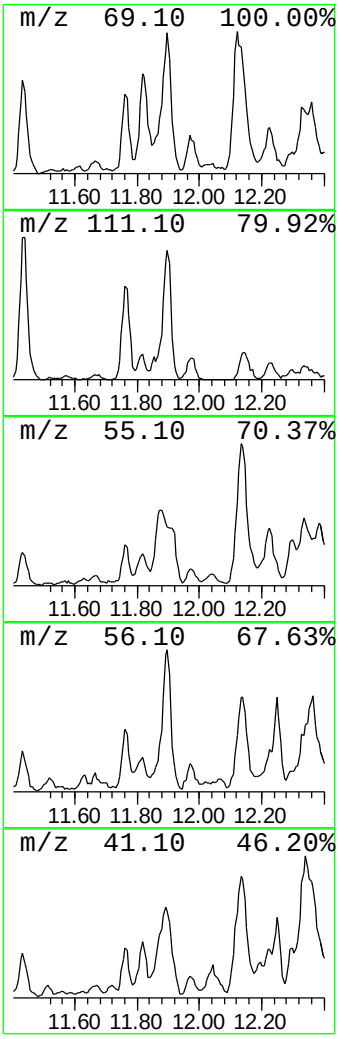
****

Peak Number 2 Cyclohexane, 1,1,2-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	26.80 ug/l	458669	Chlorobenzene-d5	11.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	90
2	Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	74
3	Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	58
4	Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	52
5	trans-2-Methyl-3-octene	126	C9H18	052937-36-7	52





Library Search Compound Report

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Misc : 7.10G/5ML/MSVOA W/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : TP-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W102319S.M
Quant Title : SW846 8260

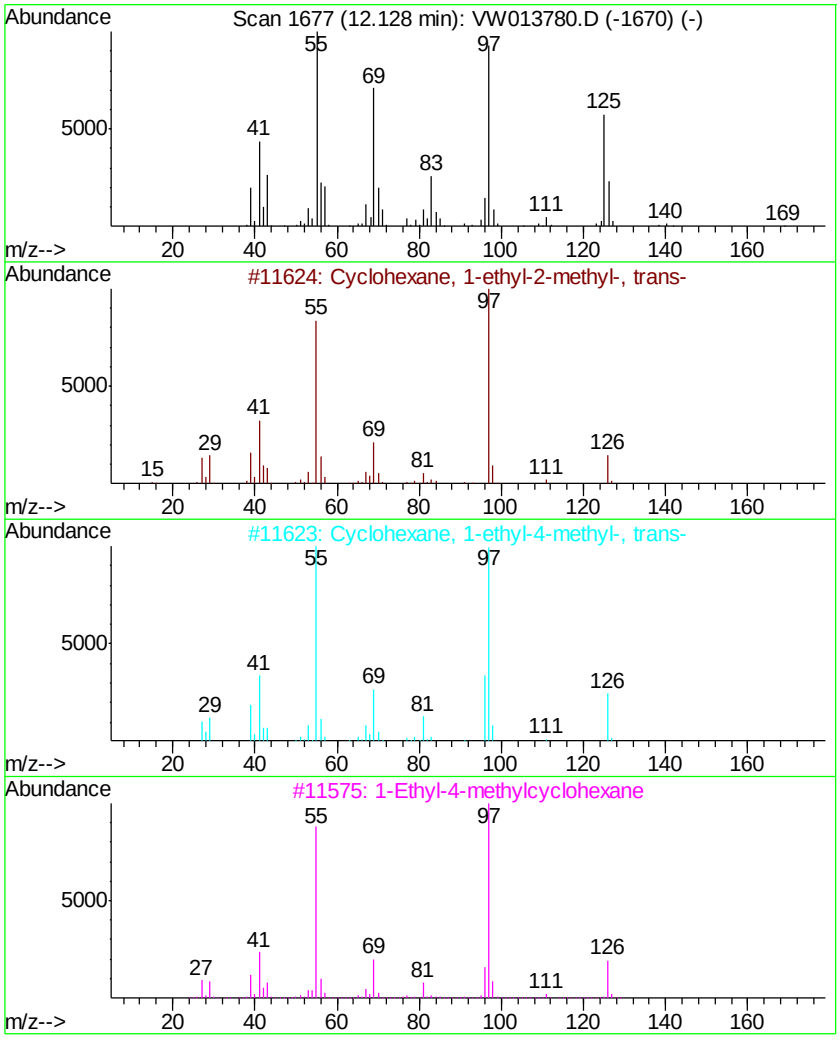
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

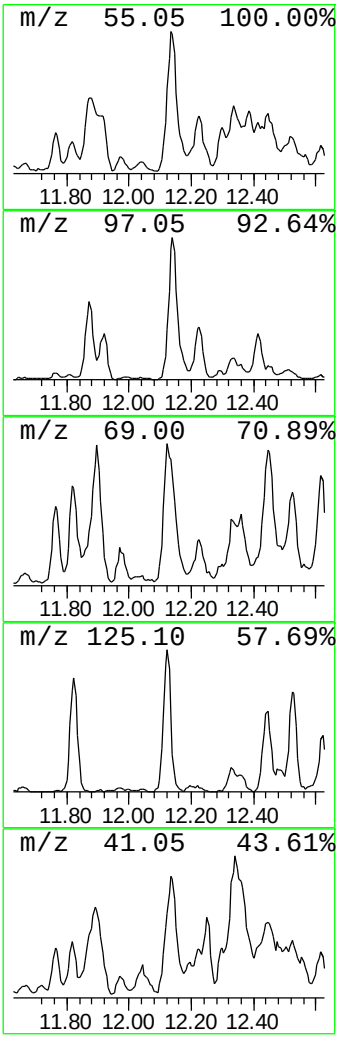
****

Peak Number 3 Cyclohexane, 1-ethyl-2-meth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	31.70 ug/l	542547	Chlorobenzene-d5	11.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	62
2	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	58
3	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	58
4	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	58
5	3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	58





Library Search Compound Report

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ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : TP-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W102319S.M
Quant Title : SW846 8260

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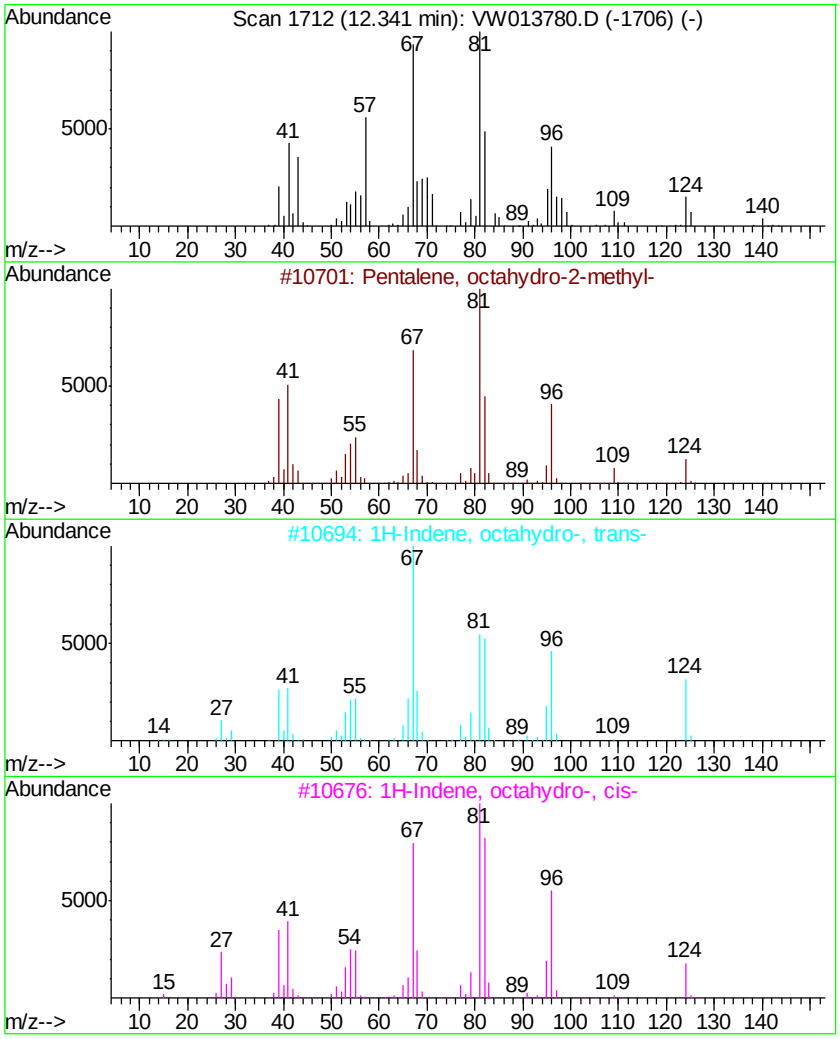
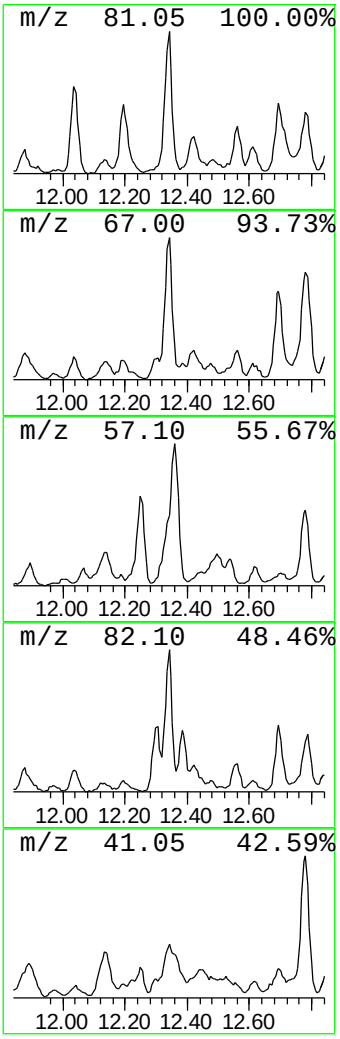
****

Peak Number 4 Pentalene, octahydro-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.34	31.42 ug/l	537740	Chlorobenzene-d5	11.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentalene, octahydro-2-methyl-	124	C9H16	003868-64-2	83
2	1H-Indene, octahydro-, trans-	124	C9H16	003296-50-2	58
3	1H-Indene, octahydro-, cis-	124	C9H16	004551-51-3	52
4	Cyclopropane, (2-methylenebutyl)-	110	C8H14	074685-56-6	52
5	Spiro[2.5]octane-1,1-dicarbonitr...	174	C11H14N2	1000151-43-1	43





Library Search Compound Report

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ALS Vial : 23 Sample Multiplier: 1

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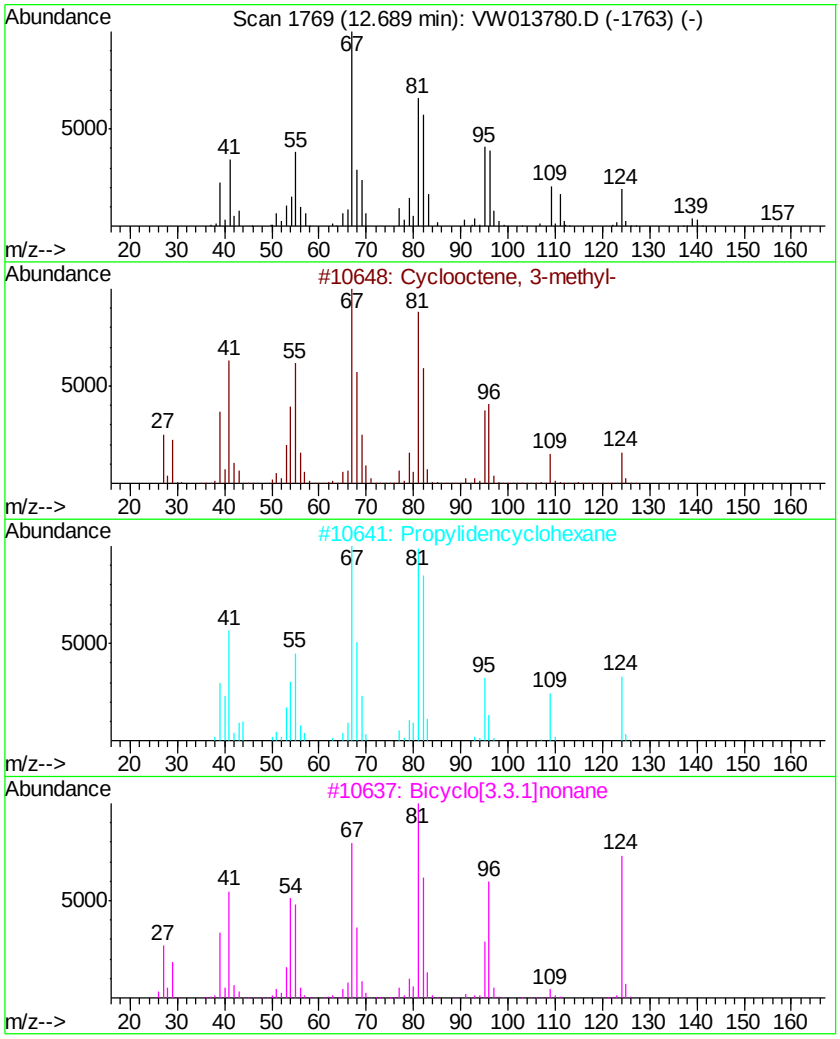
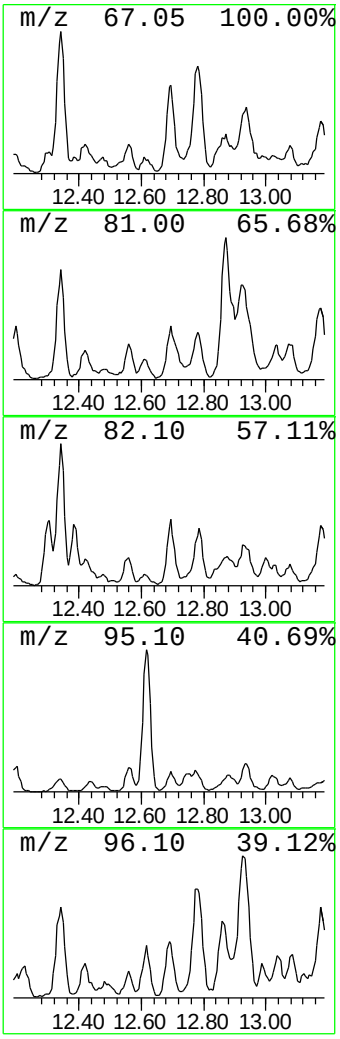
****

Peak Number 5 Cyclooctene, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.69	23.75 ug/l	334620	1,4-Dichlorobenzene-d4	13.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclooctene, 3-methyl-	124	C9H16	013152-05-1	64
2	Propylidencyclohexane	124	C9H16	002129-93-3	64
3	Bicyclo[3.3.1]nonane	124	C9H16	000280-65-9	50
4	1-Methylcyclooctene	124	C9H16	000933-11-9	49
5	3-Octyne, 2-methyl-	124	C9H16	055402-15-8	49





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Quant Title : SW846 8260

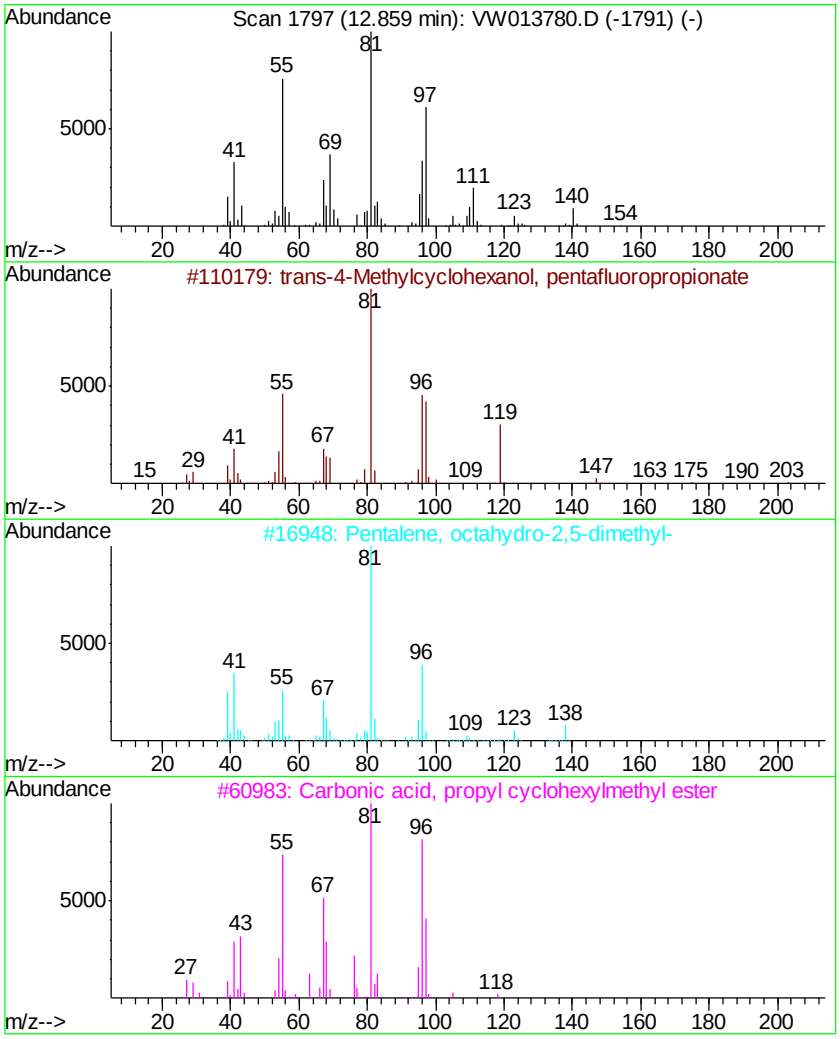
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

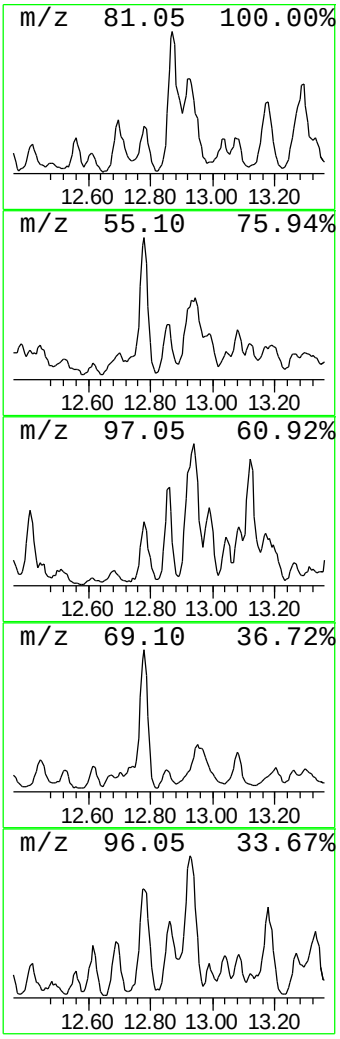
****

Peak Number 6 trans-4-Methylcyclohexanol,... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.86	30.66 ug/l	431987	1,4-Dichlorobenzene-d4	13.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-4-Methylcyclohexanol, pent...	260	C10H13F5O2	1000364-13-9	53
2	Pentalene, octahydro-2,5-dimethyl-	138	C10H18	028588-55-8	47
3	Carbonic acid, propyl cyclohexyl...	200	C11H20O3	1000357-81-6	47
4	Formic acid, cis-4-methylcyclohe...	142	C8H14O2	1000368-24-7	43
5	Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	41





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ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : TP-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W102319S.M
Quant Title : SW846 8260

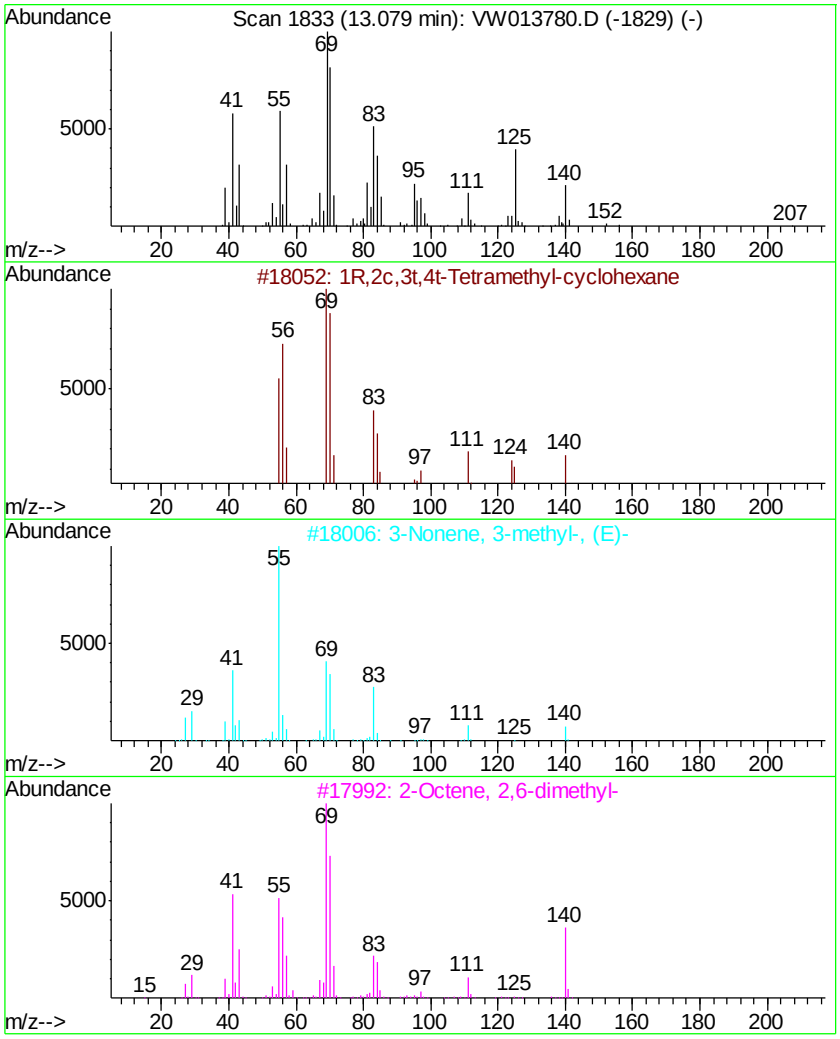
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

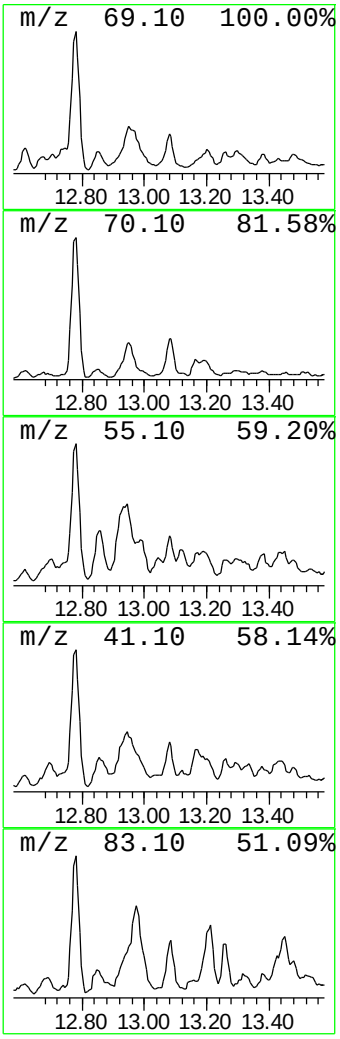
****

Peak Number 7 1R,2c,3t,4t-Tetramethyl-cyc... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.08	16.07 ug/l	226410	1,4-Dichlorobenzene-d4	13.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1R,2c,3t,4t-Tetramethyl-cyclohexane	140	C10H20	1000144-07-3	91
2	3-Nonene, 3-methyl-, (E)-	140	C10H20	069405-42-1	58
3	2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	53
4	2-Octene, 4-ethyl-	140	C10H20	053966-52-2	43
5	2-Methylene cyclopentanol	98	C6H10O	020461-31-8	38





Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW102819\
Data File : VW013780.D
Acq On : 28 Oct 2019 21:09
Operator : SY/VA
Sample : K5563-11
Misc : 7.10G/5ML/MSVOA W/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : TP-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W102319S.M
Quant Title : SW846 8260

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

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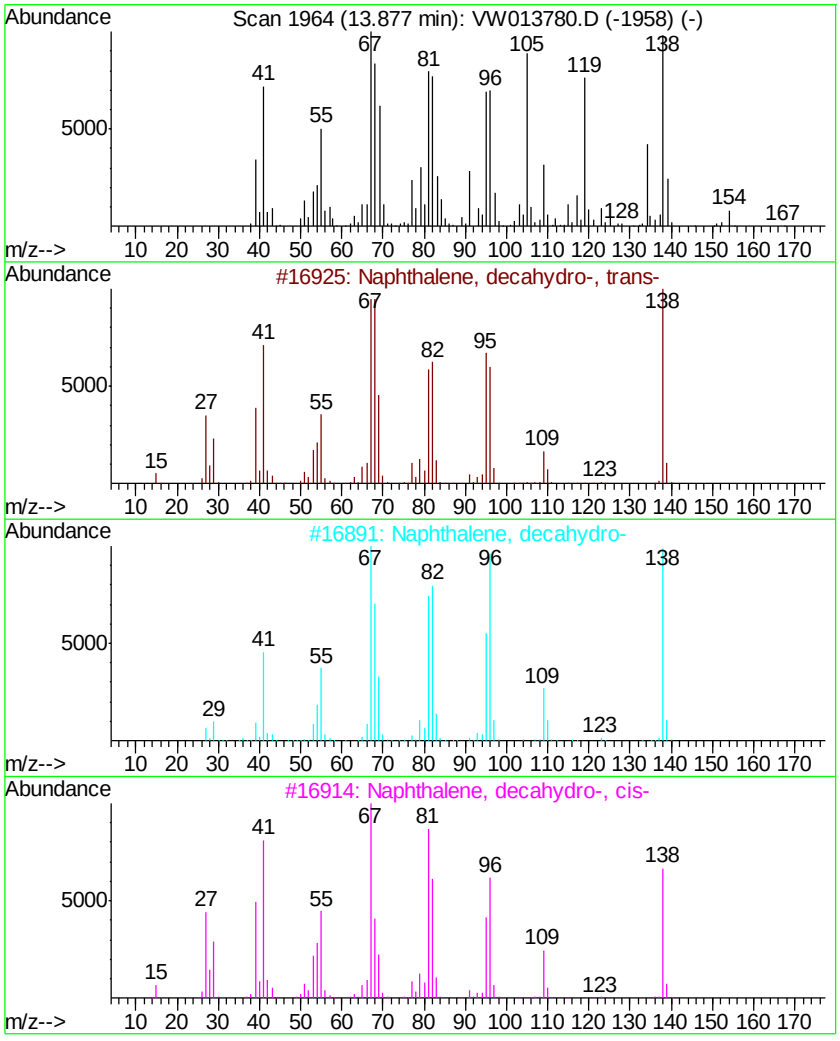
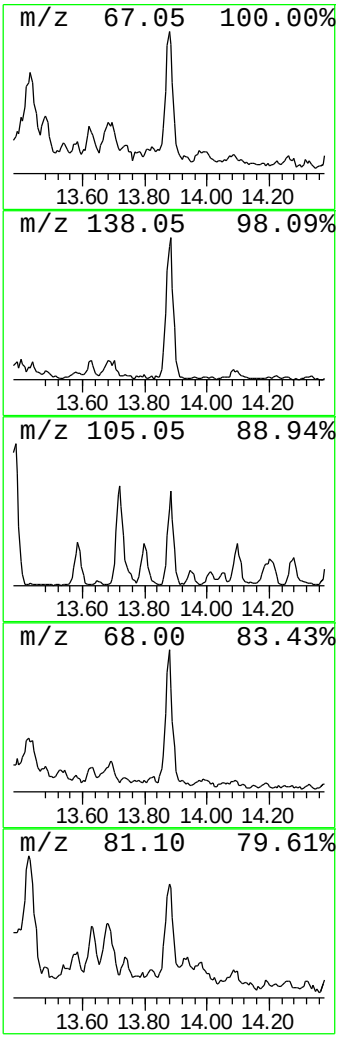
Peak Number 8 Naphthalene, decahydro-, tr... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.88	29.38 ug/l	413898	1,4-Dichlorobenzene-d4	13.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	96
2	Naphthalene, decahydro-	138	C10H18	000091-17-8	86
3	Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	64
4	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	56
5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	56





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW102819\
Data File : VW013780.D
Acq On : 28 Oct 2019 21:09
Operator : SY/VA
Sample : K5563-11
Misc : 7.10G/5ML/MSVOA W/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : TP-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W102319S.M
Quant Title : SW846 8260

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

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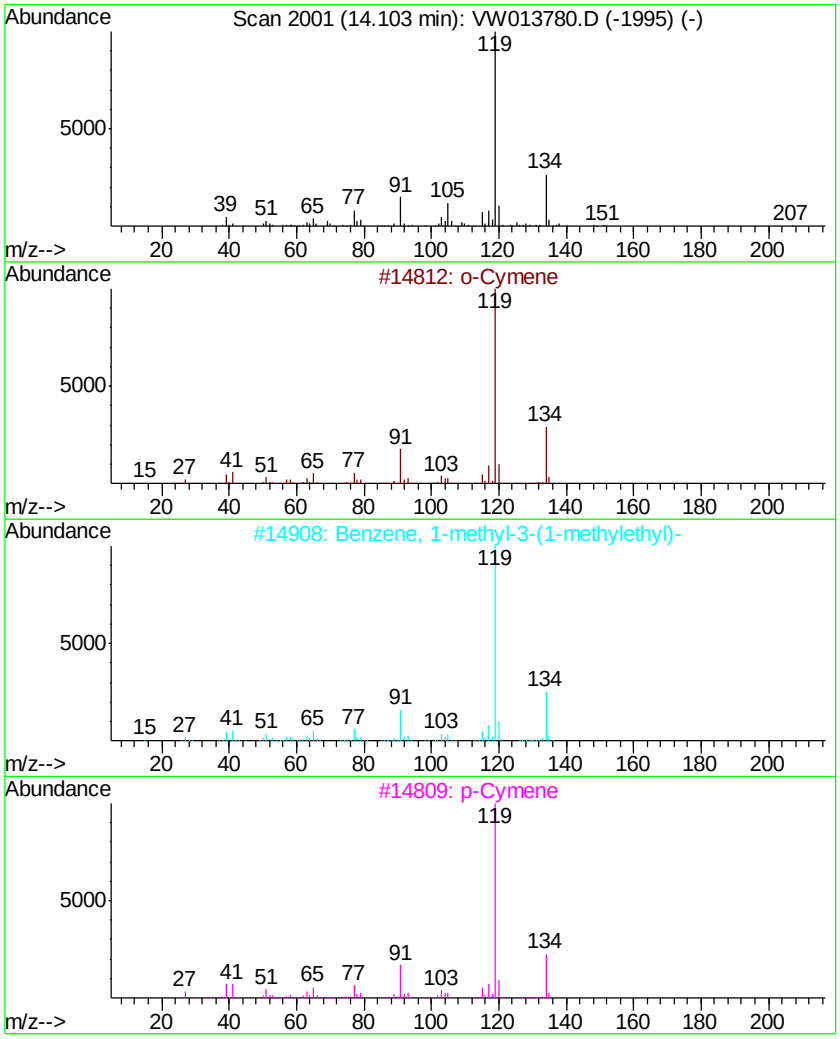
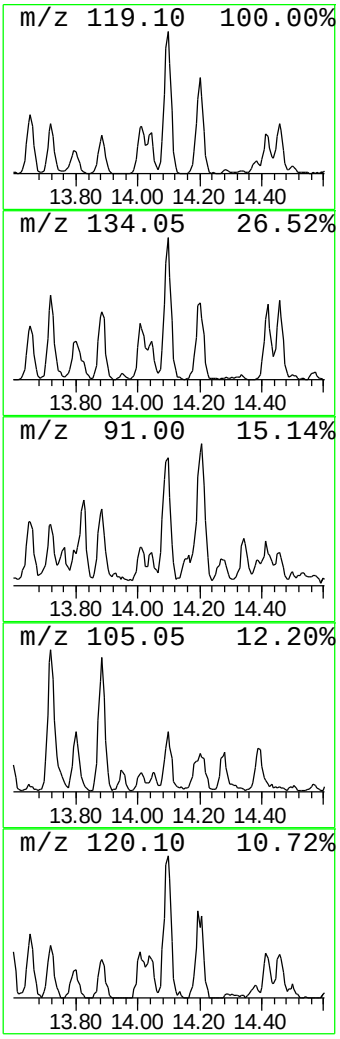
Peak Number 9 o-Cymene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.10	18.21 ug/l	256554	1,4-Dichlorobenzene-d4	13.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	95
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
3	p-Cymene	134	C10H14	000099-87-6	94
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW102819\
Data File : VW013780.D
Acq On : 28 Oct 2019 21:09
Operator : SY/VA
Sample : K5563-11
Misc : 7.10G/5ML/MSVOA W/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : TP-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W102319S.M
Quant Title : SW846 8260

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

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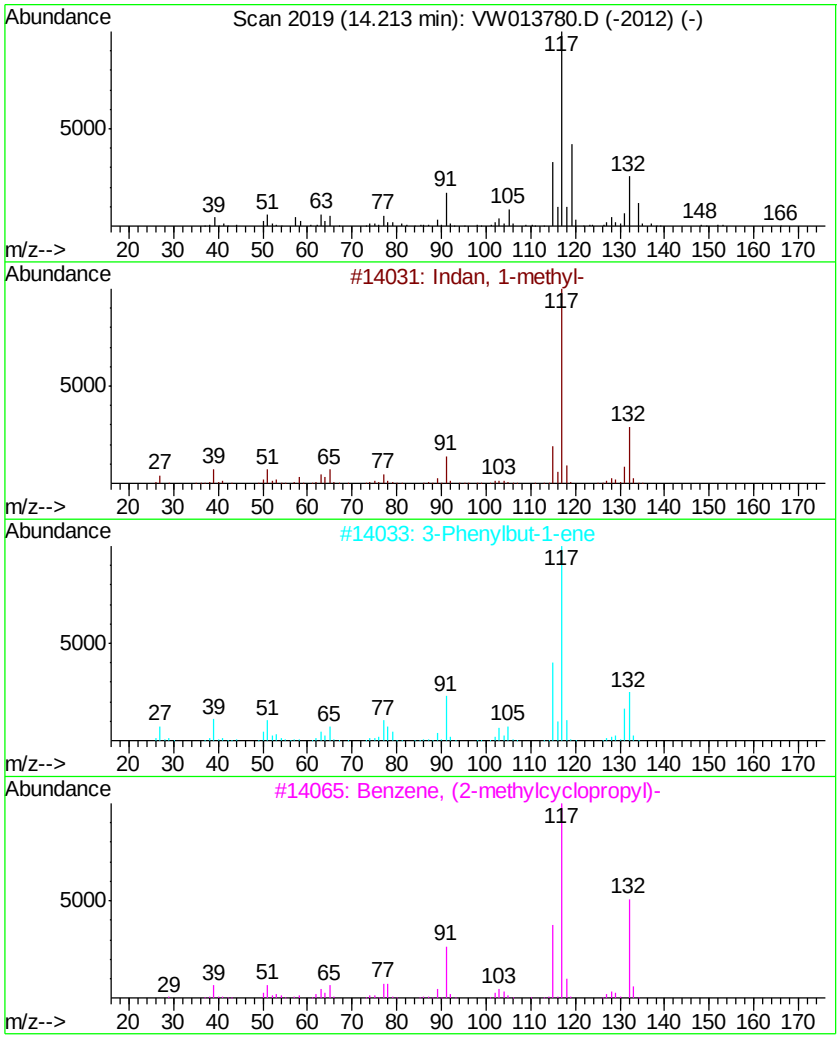
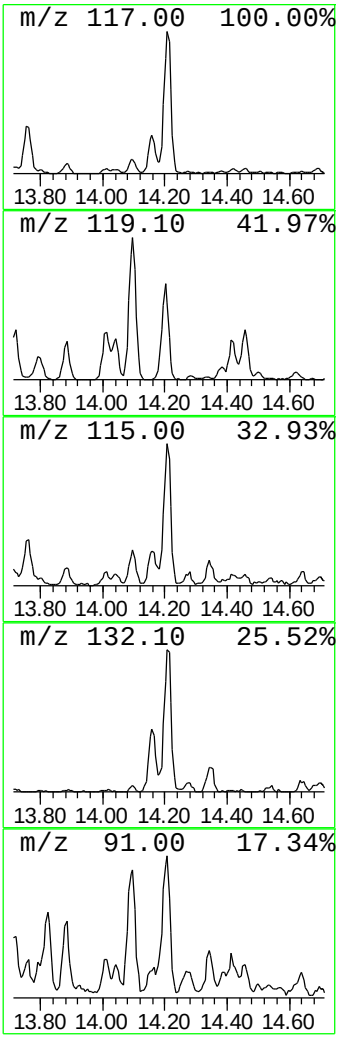
Peak Number 10 Indan, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	24.15 ug/l	340156	1,4-Dichlorobenzene-d4	13.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	70
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	64
3	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	58
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	58
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	58





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW102819\
Data File : VW013780.D
Acq On : 28 Oct 2019 21:09
Operator : SY/VA
Sample : K5563-11
Misc : 7.10G/5ML/MSVOA_W/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : TP-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W102319S.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
Cyclohexane, 1,1,...	11.23	16.3	ug/l	279204	3	11.63	855657	50.0
Cyclohexane, 1,1,...	11.90	26.8	ug/l	458669	3	11.63	855657	50.0
Cyclohexane, 1-et...	12.13	31.7	ug/l	542547	3	11.63	855657	50.0
Pentalene, octahy...	12.34	31.4	ug/l	537740	3	11.63	855657	50.0
Cyclooctene, 3-me...	12.69	23.8	ug/l	334620	4	13.55	704390	50.0
trans-4-Methylcyc...	12.86	30.7	ug/l	431987	4	13.55	704390	50.0
1R,2c,3t,4t-Tetra...	13.08	16.1	ug/l	226410	4	13.55	704390	50.0
Naphthalene, deca...	13.88	29.4	ug/l	413898	4	13.55	704390	50.0
o-Cymene	14.10	18.2	ug/l	256554	4	13.55	704390	50.0
Indan, 1-methyl-	14.21	24.1	ug/l	340156	4	13.55	704390	50.0