

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091218\
 Data File : VX004490.D
 Acq On : 12 Sep 2018 13:13
 Operator : JC/MD
 Sample : J4848-08
 Misc : 4.37g/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EP1-MW-B(16-20)

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.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	1.636	63	79	80	rBV3	13145	44988	1.04%	0.081%
2	5.507	701	714	727	rBV2	122707	352810	8.12%	0.635%
3	5.659	727	739	754	rVB	174558	494502	11.38%	0.890%
4	6.068	795	806	819	rBV	134949	356682	8.21%	0.642%
5	6.860	926	936	951	rBV	277125	656764	15.12%	1.181%
6	8.714	1225	1240	1255	rBV	676808	1108916	25.52%	1.995%
7	9.037	1287	1293	1303	rVB2	78468	132034	3.04%	0.238%
8	9.445	1349	1360	1369	rVV	125302	209691	4.83%	0.377%
9	9.555	1373	1378	1393	rVV2	118884	257019	5.92%	0.462%
10	9.683	1393	1399	1400	rVV2	31279	43974	1.01%	0.079%
11	9.713	1400	1404	1408	rVV	71706	144384	3.32%	0.260%
12	9.750	1408	1410	1416	rVV3	41203	65338	1.50%	0.118%
13	9.811	1416	1420	1425	rVV	29259	52021	1.20%	0.094%
14	9.890	1425	1433	1447	rVB	361273	630439	14.51%	1.134%
15	10.055	1455	1460	1465	rBV3	20099	43626	1.00%	0.078%
16	10.116	1465	1470	1475	rBV	672268	931959	21.45%	1.676%
17	10.171	1475	1479	1485	rVB3	56512	103963	2.39%	0.187%
18	10.238	1485	1490	1496	rBV2	196081	342220	7.88%	0.616%
19	10.335	1502	1506	1508	rBV	178870	265679	6.12%	0.478%
20	10.408	1515	1518	1530	rVB5	133742	281650	6.48%	0.507%
21	10.512	1530	1535	1541	rBV	306894	457808	10.54%	0.824%
22	10.591	1543	1548	1550	rBV2	131967	205048	4.72%	0.369%
23	10.646	1550	1557	1561	rBV2	739905	1428490	32.88%	2.570%
24	10.725	1564	1570	1572	rBV4	162504	292670	6.74%	0.526%
25	10.768	1572	1577	1580	rVB4	316114	503138	11.58%	0.905%
26	10.835	1580	1588	1592	rBV	2398290	3492891	80.40%	6.283%
27	10.914	1592	1601	1604	rBV3	866115	1709521	39.35%	3.075%
28	10.951	1604	1607	1611	rVB	1361422	1744256	40.15%	3.138%
29	11.000	1611	1615	1621	rVB2	429726	696097	16.02%	1.252%
30	11.061	1621	1625	1634	rBV5	533079	1349976	31.07%	2.428%
31	11.140	1634	1638	1642	rVB2	727453	992836	22.85%	1.786%
32	11.183	1642	1645	1649	rVB	294474	360568	8.30%	0.649%
33	11.274	1649	1660	1671	rBV2	2102418	4344455	100.00%	7.815%
34	11.396	1675	1680	1684	rBV3	830242	1624122	37.38%	2.921%

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35	11.445	1684	1688	1692	rVB3	883142	1239595	28.53%	2.230%
36	11.487	1692	1695	1698	rBV2	436687	457528	10.53%	0.823%
37	11.536	1698	1703	1707	rVB3	880098	1207817	27.80%	2.173%
38	11.579	1707	1710	1716	rVB5	355190	575764	13.25%	1.036%
39	11.658	1719	1723	1725	rBV	472708	661039	15.22%	1.189%
40	11.689	1725	1728	1732	rVB4	823891	1091777	25.13%	1.964%
41	11.737	1732	1736	1739	rVB4	339919	464899	10.70%	0.836%
42	11.804	1739	1747	1750	rBV4	789526	1536649	35.37%	2.764%
43	11.896	1757	1762	1764	rBV2	851906	1291996	29.74%	2.324%
44	11.945	1766	1770	1775	rVB5	1238753	2043960	47.05%	3.677%
45	12.036	1781	1785	1789	rVB2	618489	902039	20.76%	1.623%
46	12.085	1789	1793	1798	rBV2	944857	1456179	33.52%	2.619%
47	12.170	1803	1807	1811	rBV5	307365	526589	12.12%	0.947%
48	12.219	1812	1815	1817	rVV2	352211	446762	10.28%	0.804%
49	12.249	1817	1820	1824	rVB4	336207	535250	12.32%	0.963%
50	12.310	1824	1830	1834	rBV6	404845	836895	19.26%	1.505%
51	12.371	1835	1840	1847	rBV3	1490548	2828446	65.10%	5.088%
52	12.432	1847	1850	1855	rVV4	299339	610718	14.06%	1.099%
53	12.499	1855	1861	1862	rVV3	388548	787651	18.13%	1.417%
54	12.530	1862	1866	1871	rVV3	735350	1421500	32.72%	2.557%
55	12.615	1874	1880	1890	rVB6	662040	2165232	49.84%	3.895%
56	12.688	1890	1892	1896	rVB3	139586	157193	3.62%	0.283%
57	12.737	1896	1900	1904	rVB4	319520	412919	9.50%	0.743%
58	12.786	1904	1908	1911	rBV4	250960	399766	9.20%	0.719%
59	12.823	1911	1914	1919	rVV5	263505	425573	9.80%	0.766%
60	12.877	1919	1923	1929	rVB4	964064	1471075	33.86%	2.646%
61	12.957	1929	1936	1943	rBV7	244913	703760	16.20%	1.266%
62	13.042	1947	1950	1954	rVV2	652024	744421	17.13%	1.339%
63	13.097	1954	1959	1963	rVB4	163181	348198	8.01%	0.626%
64	13.152	1963	1968	1973	rBV3	187954	322175	7.42%	0.580%
65	13.194	1973	1975	1981	rVB4	99720	157638	3.63%	0.284%
66	13.261	1981	1986	1989	rBV3	109727	176702	4.07%	0.318%
67	13.304	1989	1993	1997	rVB5	96708	149794	3.45%	0.269%
68	13.353	1997	2001	2006	rVB3	205001	307894	7.09%	0.554%
69	13.402	2006	2009	2011	rBV3	50159	58907	1.36%	0.106%
70	13.438	2011	2015	2018	rBV2	155446	212411	4.89%	0.382%
71	13.475	2018	2021	2025	rVB4	106432	118751	2.73%	0.214%

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Integration Parameters: RTEINT.P

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Stop Thrs : 0 Peak Location: TOP

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72	13.639	2044	2048	2051	rBV2	83708	120497	2.77%	0.217%
73	13.682	2051	2055	2061	rVB2	81908	167592	3.86%	0.301%
74	13.761	2066	2068	2071	rVV3	40518	44715	1.03%	0.080%
75	13.810	2072	2076	2081	rVV	67333	89644	2.06%	0.161%
76	13.859	2081	2084	2087	rVB2	48165	56992	1.31%	0.103%
77	13.902	2087	2091	2098	rVV2	61802	69981	1.61%	0.126%
78	13.963	2098	2101	2109	rVB8	32659	67169	1.55%	0.121%

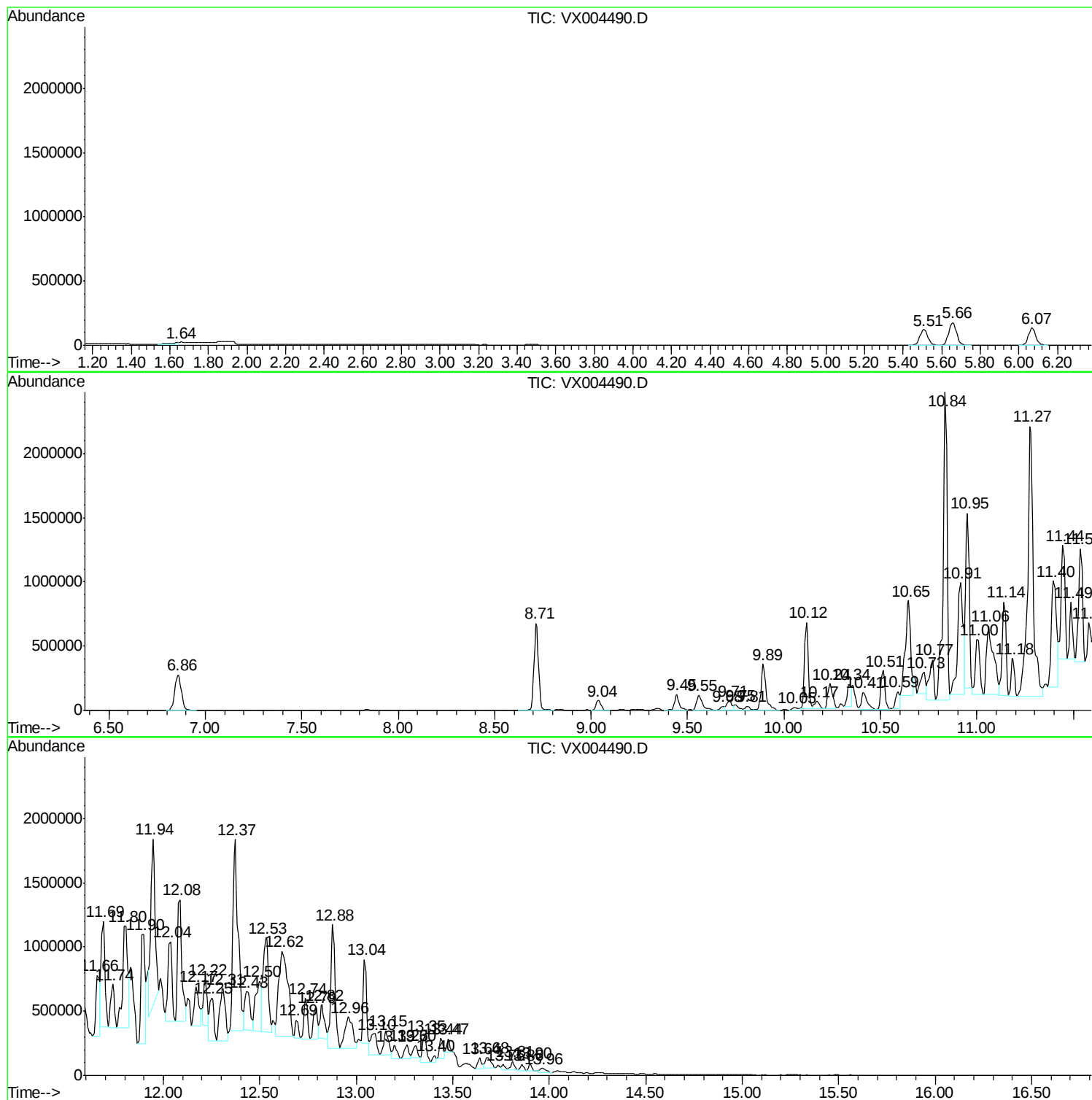
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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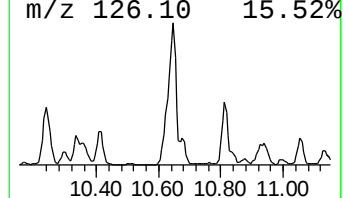
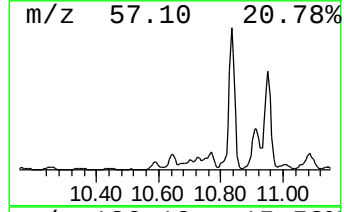
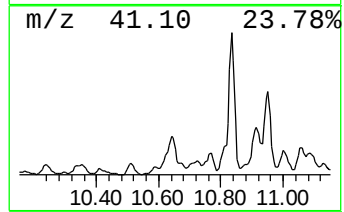
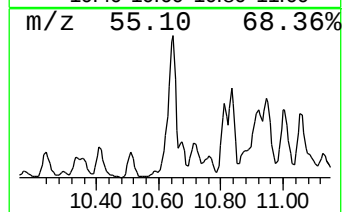
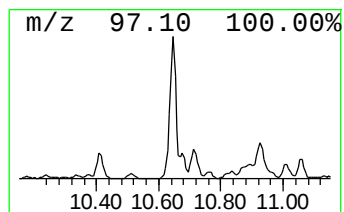
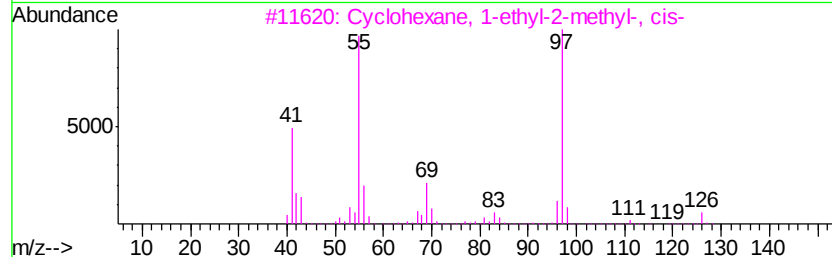
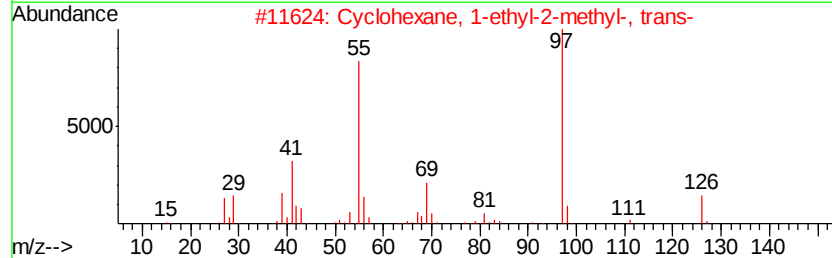
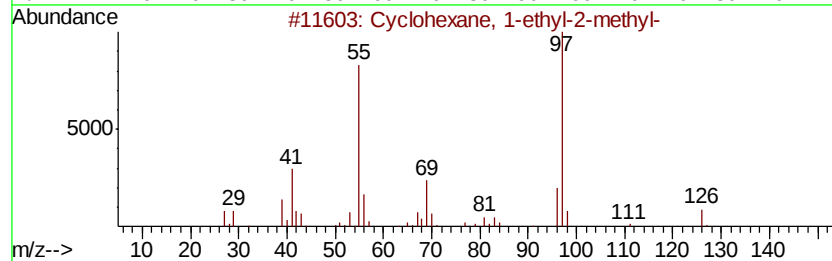
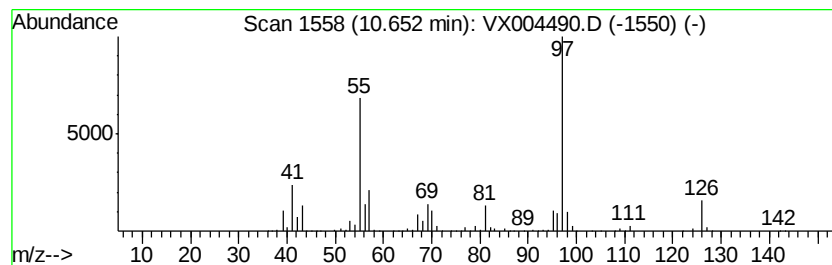
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	76.64 ug/l	1428490	Chlorobenzene-d5	10.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	90
2		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	87
3		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	86
4		1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	81
5		1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	72



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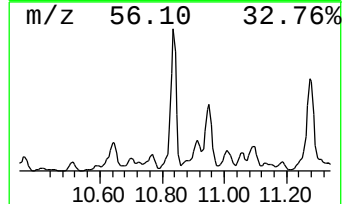
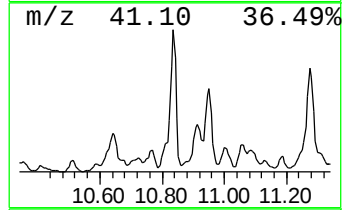
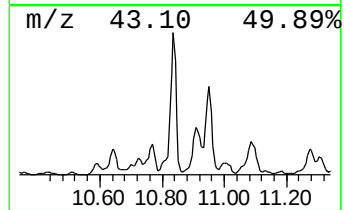
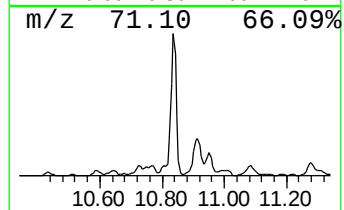
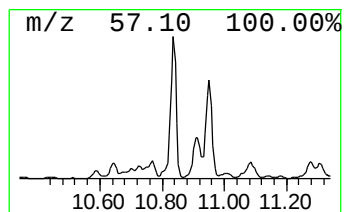
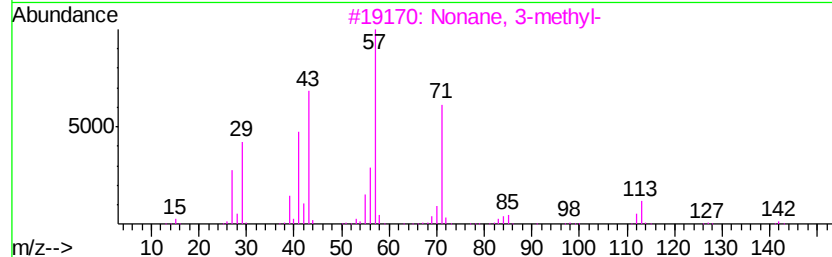
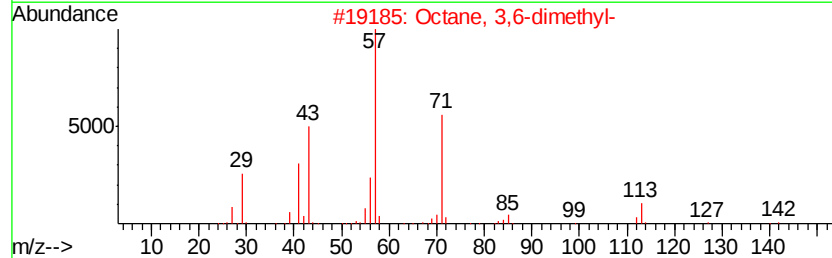
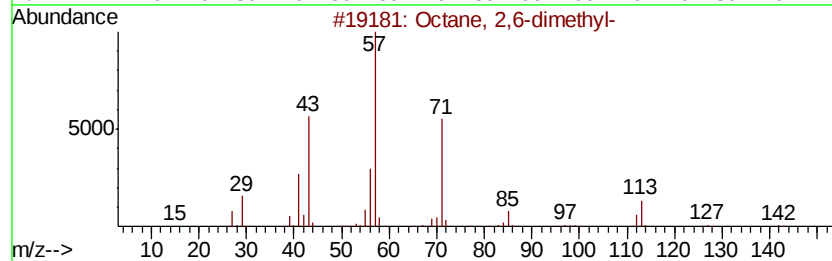
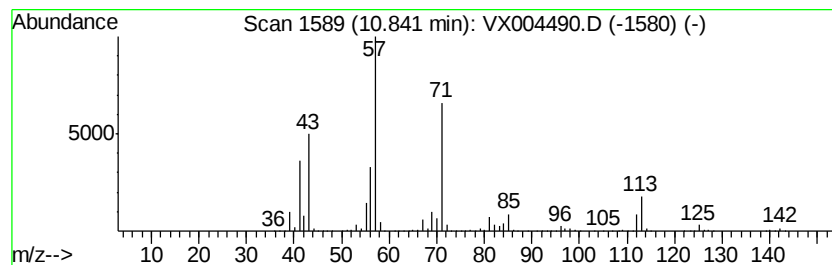
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Peak Number 2 Octane, 2,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.84	187.40 ug/l	3492890	Chlorobenzene-d5	10.12		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	94
2		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	91
3		Nonane, 3-methyl-	142	C10H22	005911-04-6	90
4		Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	53
5		Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	53



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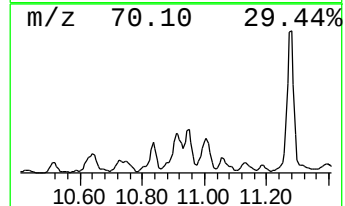
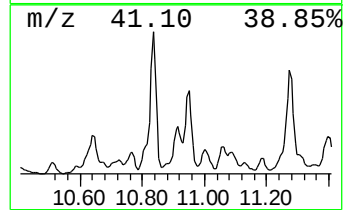
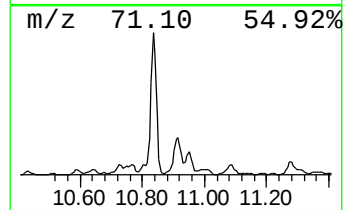
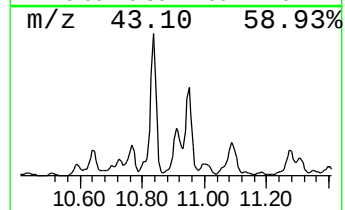
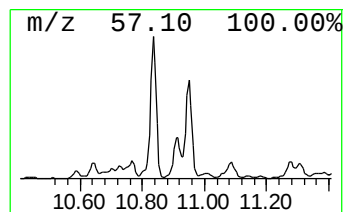
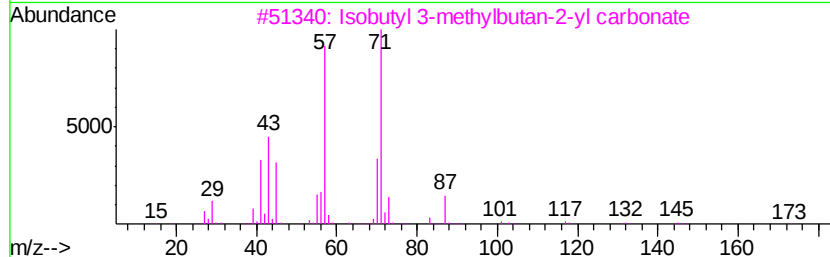
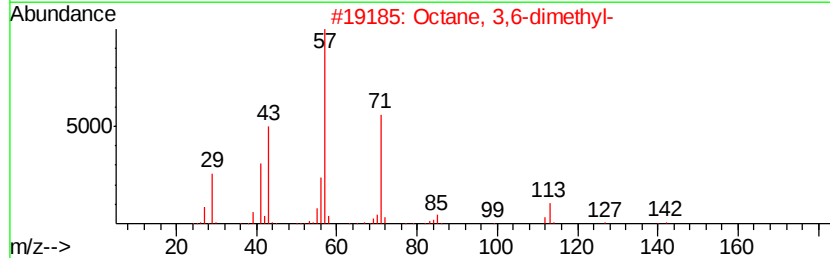
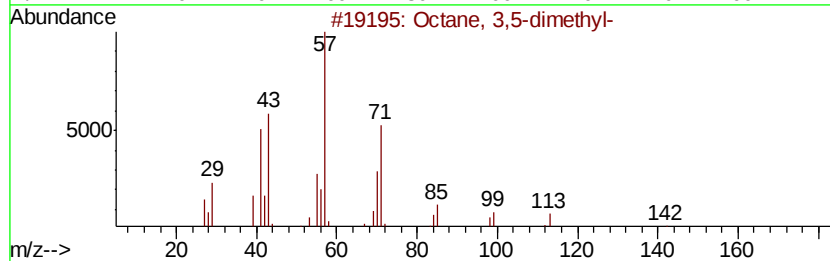
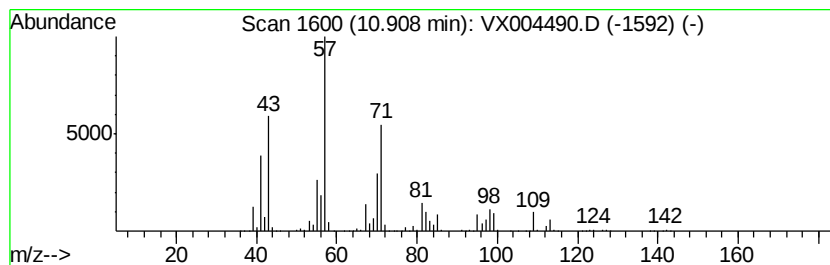
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Peak Number 3 Octane, 3,5-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.91	91.72 ug/l	1709520	Chlorobenzene-d5	10.12
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Octane, 3,5-dimethyl-	142 C10H22	015869-93-9 52
2		Octane, 3,6-dimethyl-	142 C10H22	015869-94-0 49
3		Isobutyl 3-methylbutan-2-yl carb...	188 C10H20O3	1000372-77-0 47
4		Butane, 2,2-dimethyl-	86 C6H14	000075-83-2 38
5		2,4,4-Trimethyl-1-hexene	126 C9H18	051174-12-0 38



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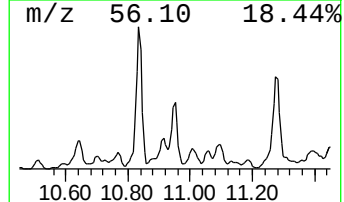
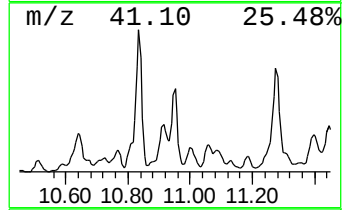
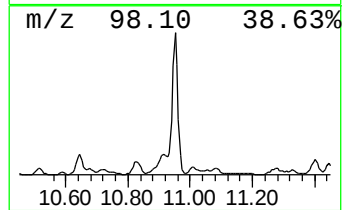
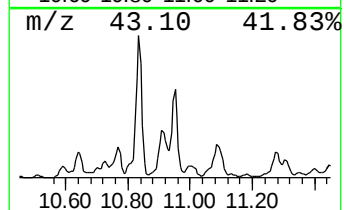
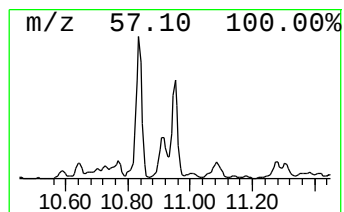
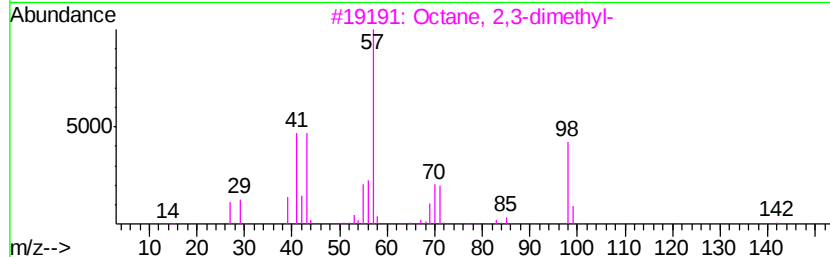
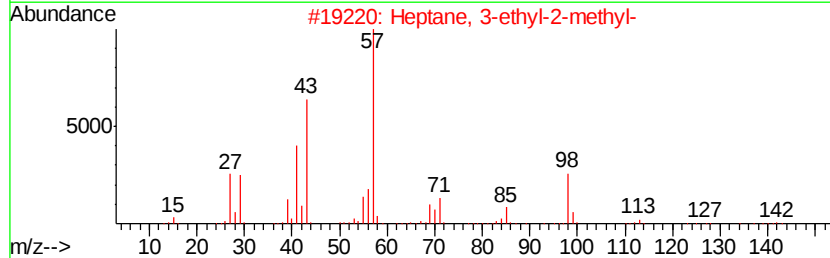
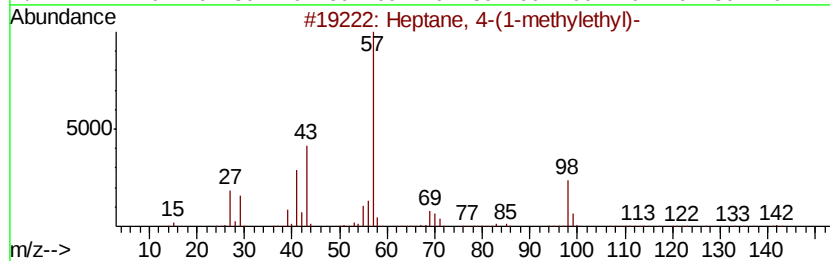
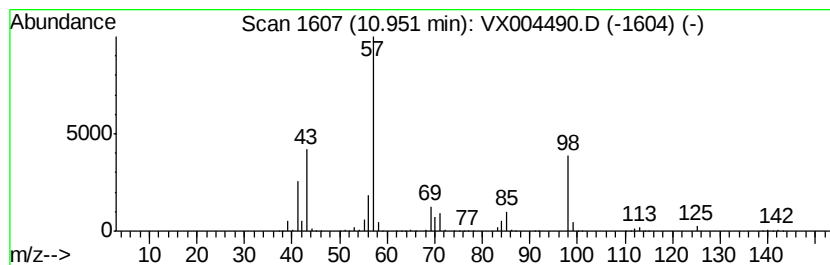
Instrument :
MSVOA_X
ClientSampled :
EP1-MW-B(16-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 4 Heptane, 4-(1-methylethyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.95	93.58 ug/l	1744260	Chlorobenzene-d5	10.12		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	64
2		Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	53
3		Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	53
4		Octane, 3-methyl-	128	C9H20	002216-33-3	43
5		Sulfurous acid, 2-ethylhexyl hep...	432	C25H52O3S	1000309-20-0	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091218\
Data File : VX004490.D
Acq On : 12 Sep 2018 13:13
Operator : JC/MD
Sample : J4848-08
Misc : 4.37µL/100µL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

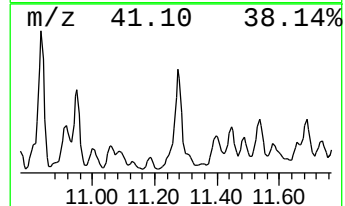
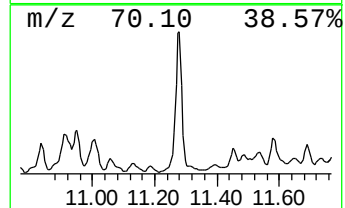
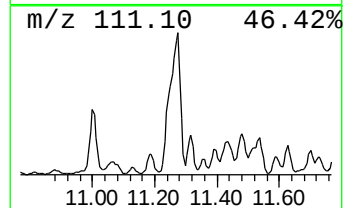
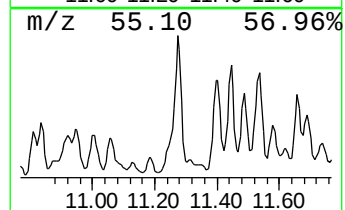
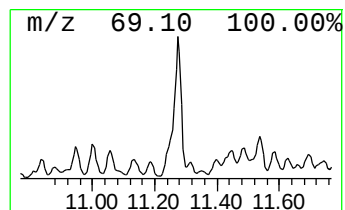
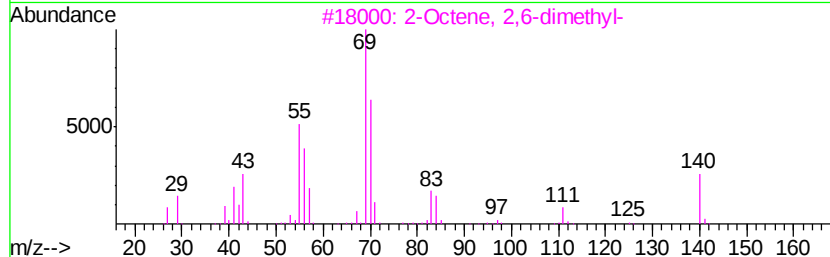
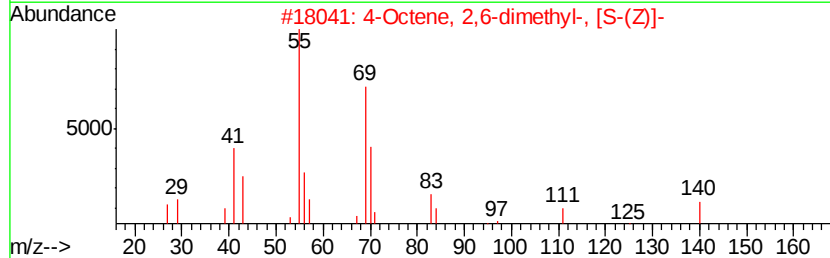
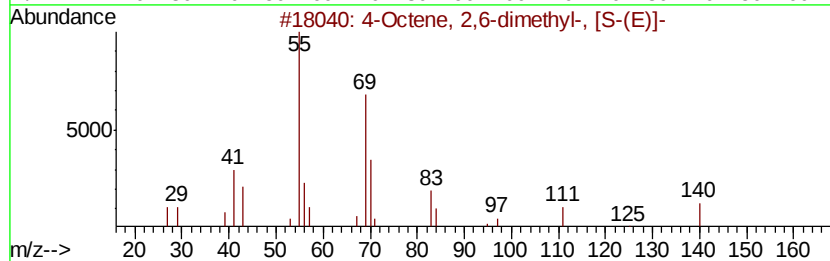
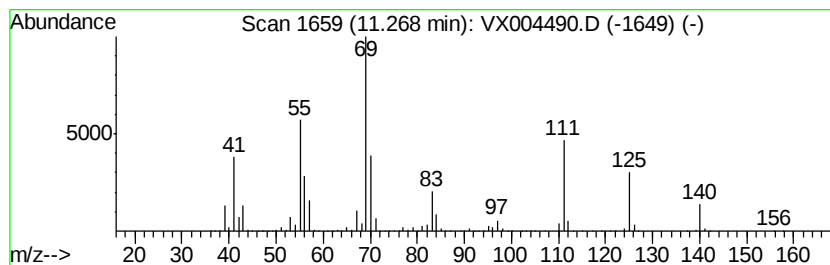
Instrument :
MSVOA_X
ClientSampleID :
EP1-MW-B(16-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 5 4-Octene, 2,6-dimethyl-, [S... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.27	149.17 ug/l	4344460	1,4-Dichlorobenzene-d4	12.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Octene, 2,6-dimethyl-, [S-(E)]-	140	C10H20	062960-76-3	76
2	4-Octene, 2,6-dimethyl-, [S-(Z)]-	140	C10H20	062960-77-4	76
3	2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	70
4	Cyclohexane, 1,2,3-trimethyl-, (...)	126	C9H18	001678-81-5	62
5	1-Hexene, 3,3,5-trimethyl-	126	C9H18	013427-43-5	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091218\
Data File : VX004490.D
Acq On : 12 Sep 2018 13:13
Operator : JC/MD
Sample : J4848-08
Misc : 4.37µ/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

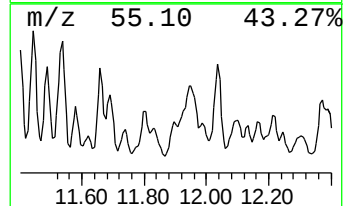
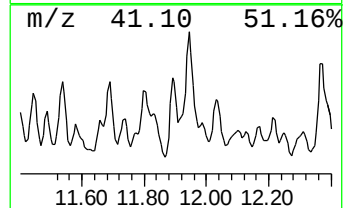
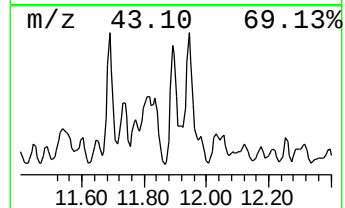
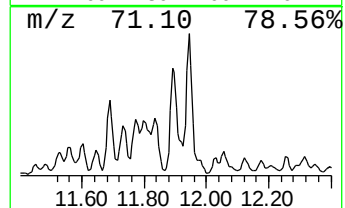
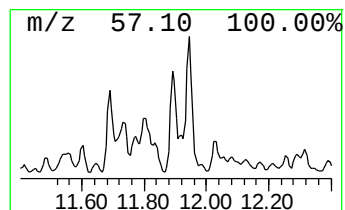
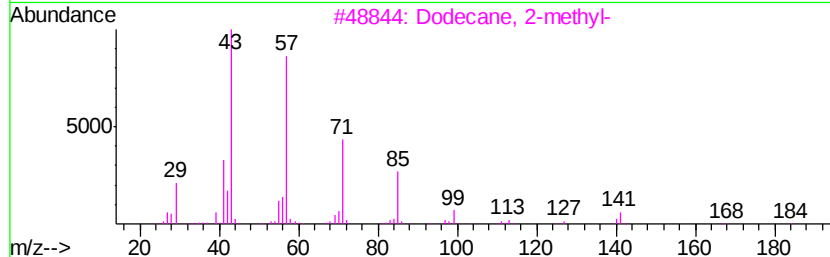
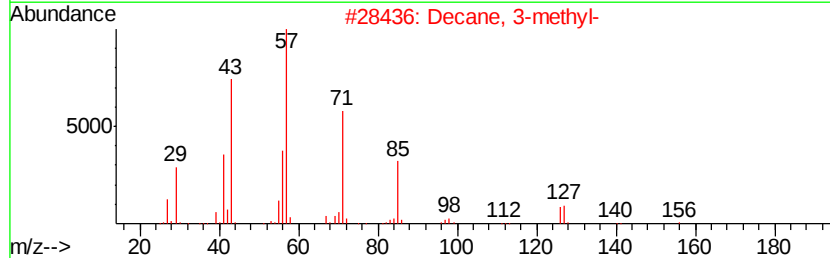
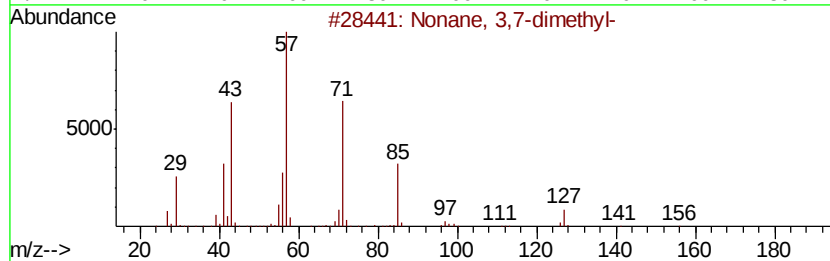
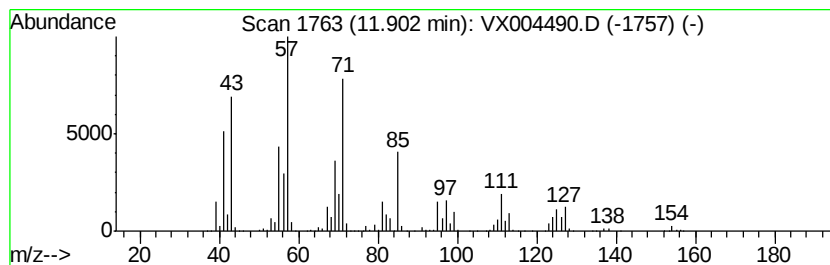
Instrument :
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ClientSampleId :
EP1-MW-B(16-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 6 Nonane, 3,7-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	44.36 ug/l	1292000	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Nonane, 3,7-dimethyl-	156 C11H24	017302-32-8	81
2	Decane, 3-methyl-	156 C11H24	013151-34-3	68
3	Dodecane, 2-methyl-	184 C13H28	001560-97-0	59
4	Octane, 2,3,3-trimethyl-	156 C11H24	062016-30-2	59
5	Tetradecane, 1-iodo-	324 C14H29I	019218-94-1	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091218\
Data File : VX004490.D
Acq On : 12 Sep 2018 13:13
Operator : JC/MD
Sample : J4848-08
Misc : 4.37µ/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

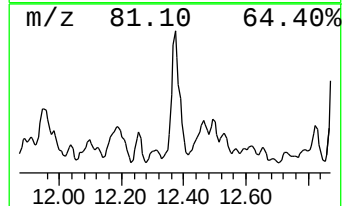
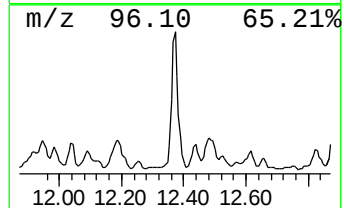
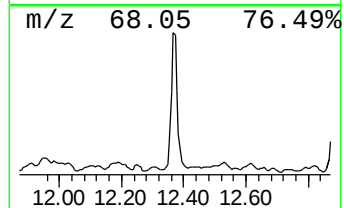
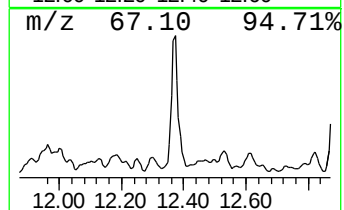
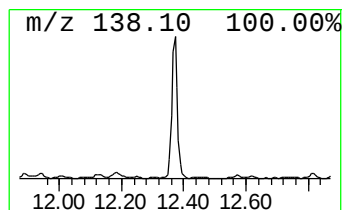
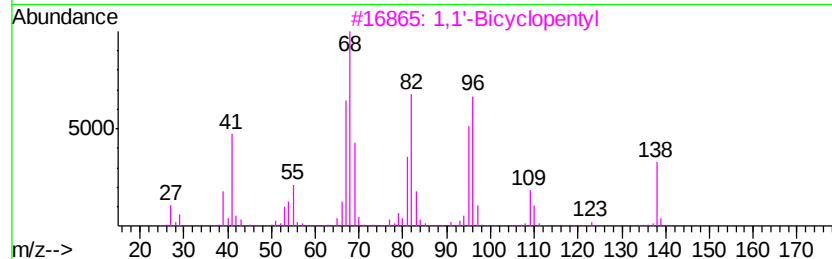
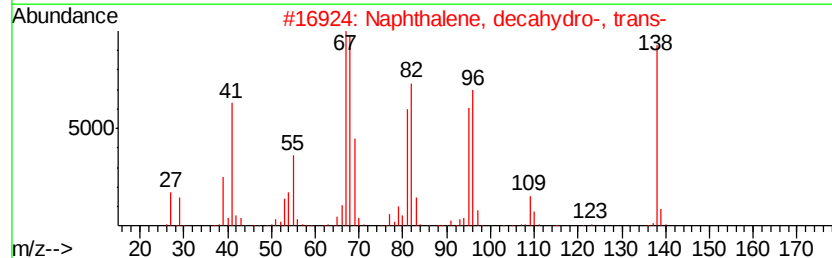
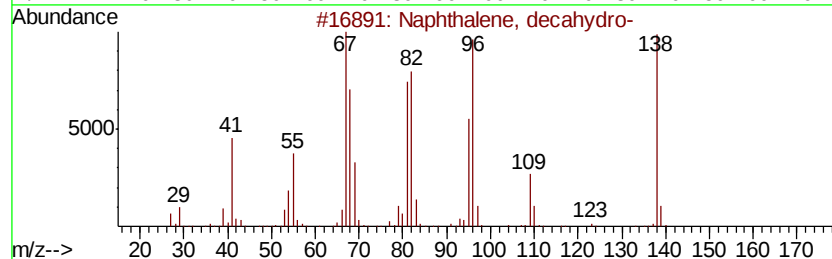
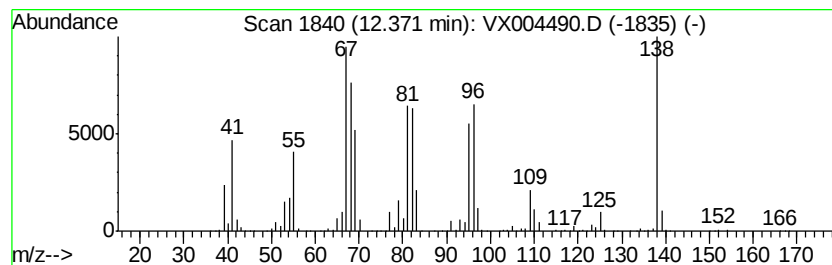
Instrument :
MSVOA_X
ClientSampleId :
EP1-MW-B(16-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, decahydro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.37	97.12 ug/l	2828450	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, decahydro-	138 C10H18	000091-17-8 97
2		Naphthalene, decahydro-, trans-	138 C10H18	000493-02-7 96
3		1,1'-Bicyclopentyl	138 C10H18	001636-39-1 95
4		Spiro[4.5]decane	138 C10H18	000176-63-6 89
5		2H-Inden-2-one, octahydro-, cis-	138 C9H14O	005689-04-3 86



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091218\
Data File : VX004490.D
Acq On : 12 Sep 2018 13:13
Operator : JC/MD
Sample : J4848-08
Misc : 4.37uL/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EP1-MW-B(16-20)

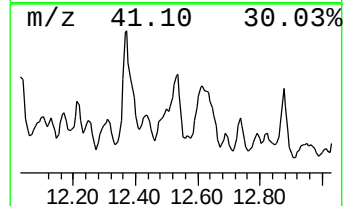
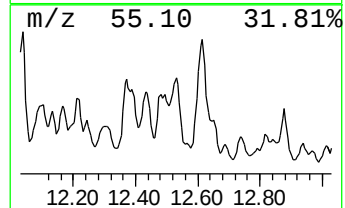
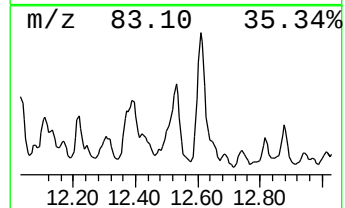
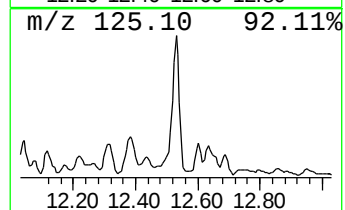
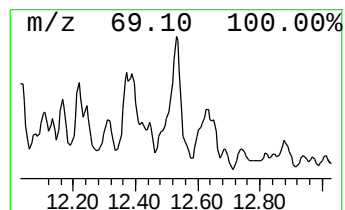
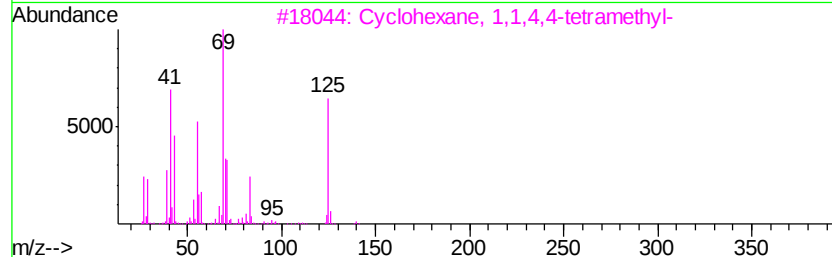
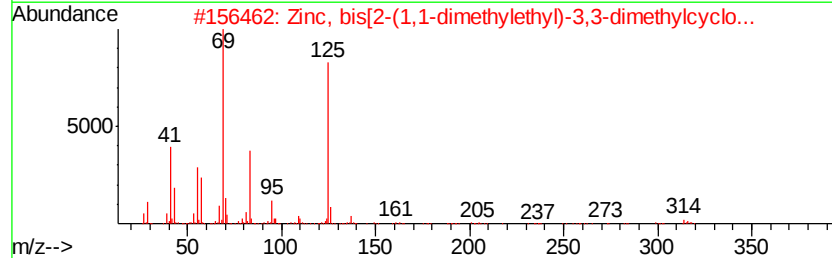
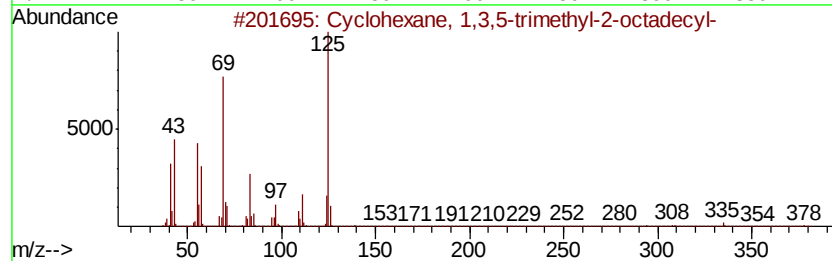
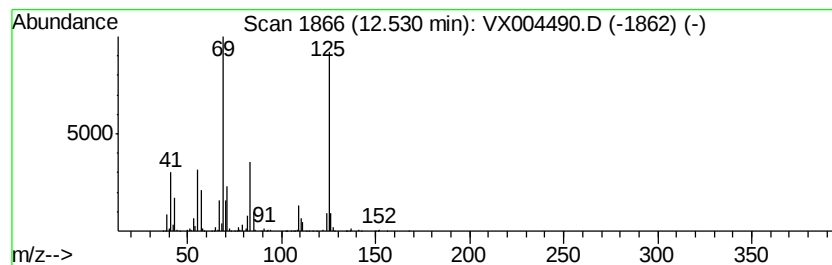
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 8 Cyclohexane, 1,3,5-trimethy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.53	48.81 ug/l	1421500	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,3,5-trimethyl-2-o...	378	C27H54	055282-34-3	64
2		Zinc, bis[2-(1,1-dimethylethyl)-...	314	C18H34Zn	074793-36-5	50
3		Cyclohexane, 1,1,4,4-tetramethyl-	140	C10H20	002223-52-1	50
4		4-Nonene, 2,3,3-trimethyl-, (Z)-	168	C12H24	063830-68-2	42
5		Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-32-9	33



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091218\
Data File : VX004490.D
Acq On : 12 Sep 2018 13:13
Operator : JC/MD
Sample : J4848-08
Misc : 4.37µ/5mL/100µL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

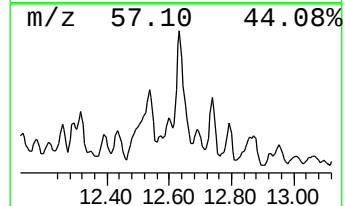
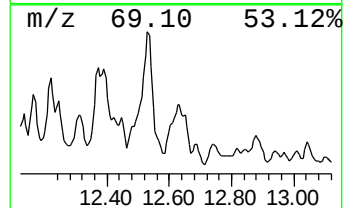
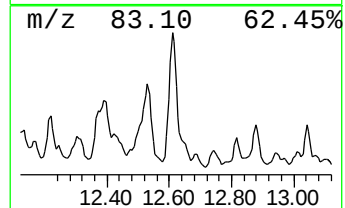
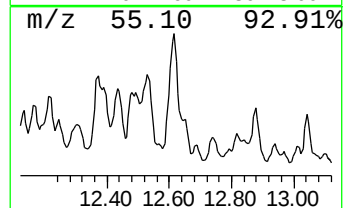
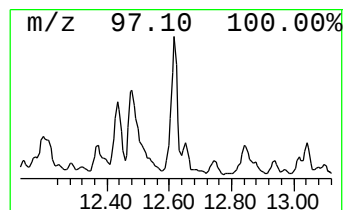
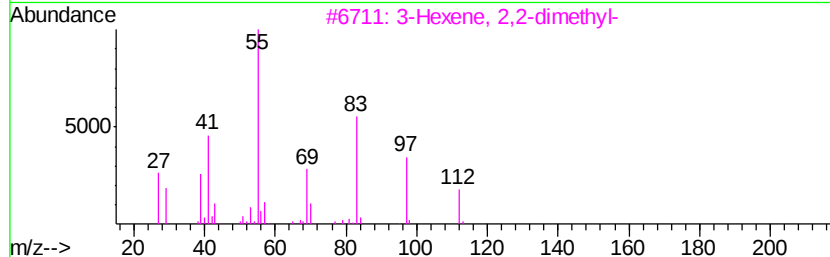
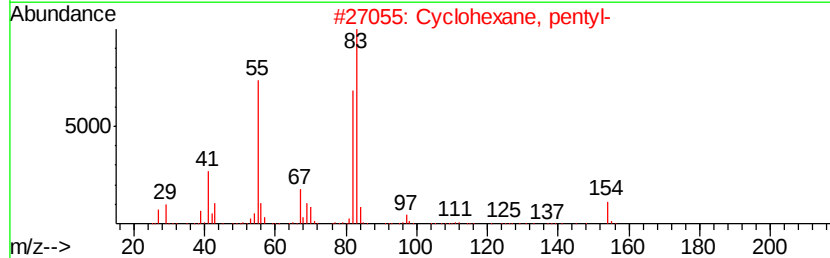
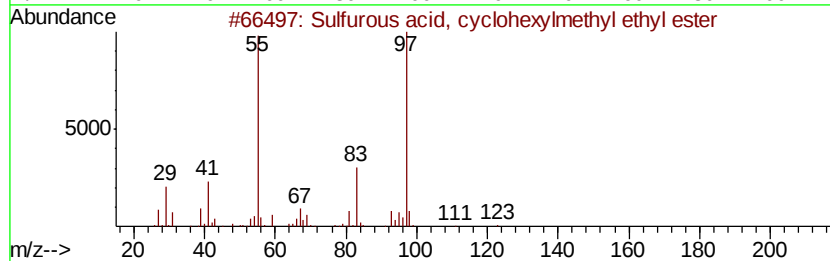
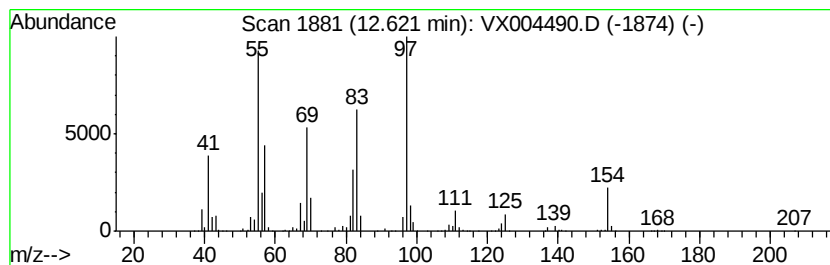
Instrument :
MSVOA_X
ClientSampleId :
EP1-MW-B(16-20)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 9 Sulfurous acid, cyclohexylm... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.62	74.35 ug/l	2165230	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Sulfurous acid, cvclohexvlmethvl...	206 C9H18O3S	1000309-21-2 59
2		Cyclohexane, pentyl-	154 C11H22	004292-92-6 53
3		3-Hexene, 2,2-dimethyl-	112 C8H16	003123-93-1 38
4		Cyclohexane, hexyl-	168 C12H24	004292-75-5 38
5		Cyclohexane, (bromomethyl)-	176 C7H13Br	002550-36-9 35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091218\
Data File : VX004490.D
Acq On : 12 Sep 2018 13:13
Operator : JC/MD
Sample : J4848-08
Misc : 4.37µ/5mL/100µL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EP1-MW-B(16-20)

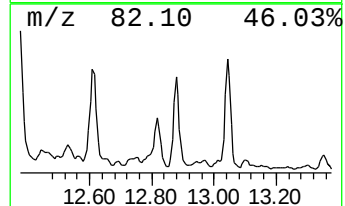
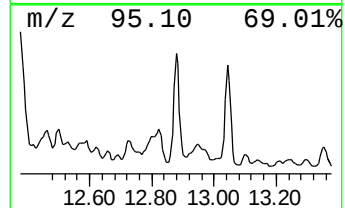
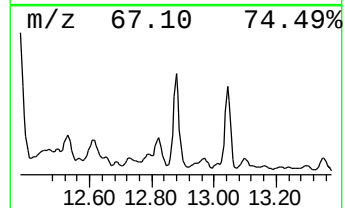
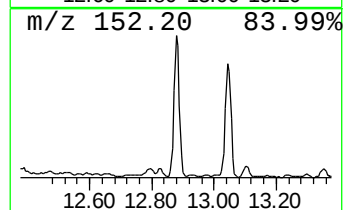
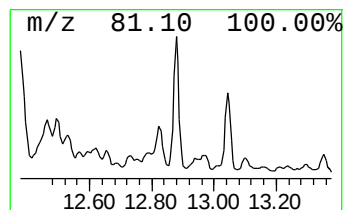
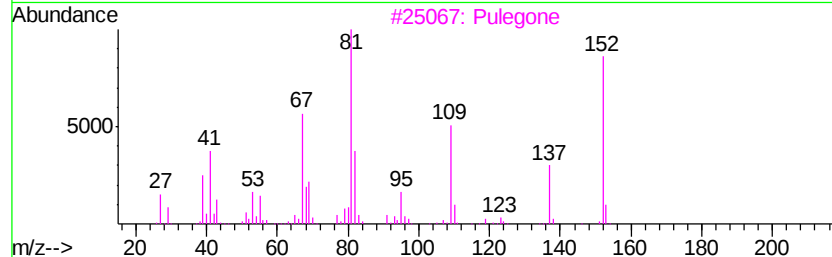
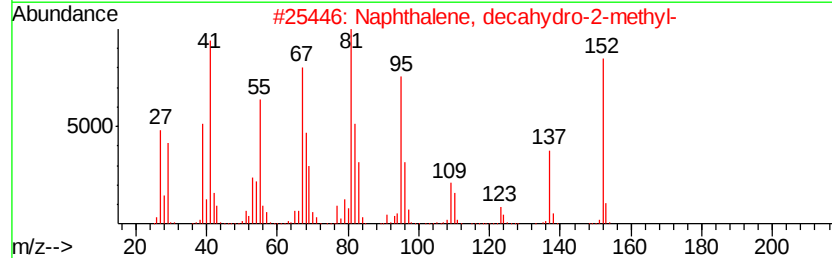
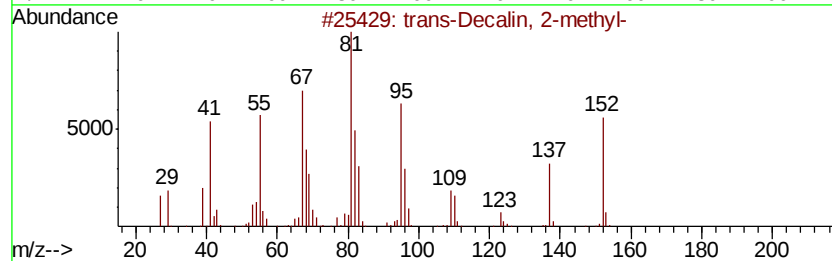
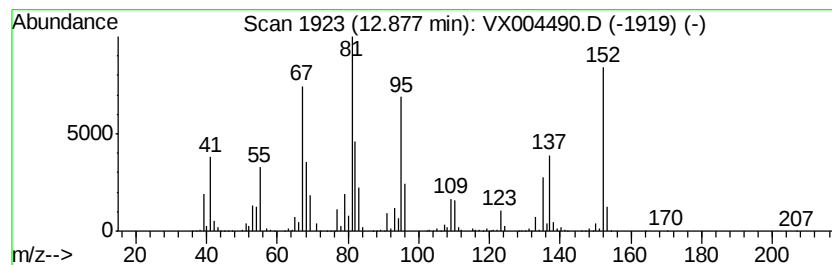
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 10 trans-Decalin, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.88	50.51 ug/l	1471080	1,4-Dichlorobenzene-d4	12.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	96
2			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	95
3			Pulegone	152	C10H16O	000089-82-7	81
4			Cyclohexanone, 5-methyl-2-(1-met...	152	C10H16O	015932-80-6	74
5			trans-4a-Methyl-decahydronaphtha...	152	C11H20	002547-27-5	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091218\
 Data File : VX004490.D
 Acq On : 12 Sep 2018 13:13
 Operator : JC/MD
 Sample : J4848-08
 Misc : 4.37g/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EP1-MW-B(16-20)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
 Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclohexane, 1-et...	10.65	76.6	ug/l	1428490	3	10.12	931959	50.0
Octane, 2,6-dimet...	10.84	187.4	ug/l	3492890	3	10.12	931959	50.0
Octane, 3,5-dimet...	10.91	91.7	ug/l	1709520	3	10.12	931959	50.0
Heptane, 4-(1-met...	10.95	93.6	ug/l	1744260	3	10.12	931959	50.0
4-Octene, 2,6-dim...	11.27	149.2	ug/l	4344460	4	12.08	1456180	50.0
Nonane, 3,7-dimet...	11.90	44.4	ug/l	1292000	4	12.08	1456180	50.0
Naphthalene, deca...	12.37	97.1	ug/l	2828450	4	12.08	1456180	50.0
Cyclohexane, 1,3,...	12.53	48.8	ug/l	1421500	4	12.08	1456180	50.0
Sulfurous acid, c...	12.62	74.3	ug/l	2165230	4	12.08	1456180	50.0
trans-Decalin, 2-...	12.88	50.5	ug/l	1471080	4	12.08	1456180	50.0