

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091218\
 Data File : VX004489.D
 Acq On : 12 Sep 2018 12:46
 Operator : JC/MD
 Sample : J4848-02
 Misc : 2.00µ/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EP1-MW-A(8-12)

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	6.860	926	936	957	rBV	271225	644417	1.14%	0.095%
2	8.665	1224	1232	1235	rBV	761895	1351807	2.39%	0.198%
3	8.713	1235	1240	1253	rVB2	779580	1649240	2.92%	0.242%
4	9.036	1287	1293	1304	rVB	734978	1231436	2.18%	0.181%
5	9.164	1304	1314	1319	rBV	356087	571134	1.01%	0.084%
6	9.347	1338	1344	1350	rVB	378294	633890	1.12%	0.093%
7	9.445	1350	1360	1369	rBV	1196816	1965589	3.48%	0.288%
8	9.561	1373	1379	1385	rBV2	1417704	2740530	4.85%	0.402%
9	9.622	1385	1389	1394	rVB	2054910	2840758	5.03%	0.417%
10	9.683	1394	1399	1403	rBV	2633360	4251039	7.53%	0.624%
11	9.817	1415	1421	1426	rBV	462565	705327	1.25%	0.103%
12	9.890	1426	1433	1438	rBV4	3302058	8071360	14.29%	1.184%
13	10.061	1456	1461	1465	rBV	1028654	1599103	2.83%	0.235%
14	10.116	1465	1470	1474	rBV2	850518	1262661	2.24%	0.185%
15	10.183	1474	1481	1486	rVB4	863723	1993299	3.53%	0.292%
16	10.244	1486	1491	1497	rBV2	2042403	3903044	6.91%	0.573%
17	10.335	1497	1506	1510	rBV2	4302145	9379490	16.61%	1.376%
18	10.384	1510	1514	1517	rVV	9519943	13127407	23.24%	1.926%
19	10.414	1517	1519	1530	rVB	7017964	11085345	19.63%	1.626%
20	10.518	1530	1536	1541	rBV	3561097	5295170	9.38%	0.777%
21	10.597	1541	1549	1550	rBV	1723252	2342782	4.15%	0.344%
22	10.652	1550	1558	1565	rBV3	17368515	30414518	53.85%	4.462%
23	10.737	1565	1572	1575	rBV4	8438001	15051215	26.65%	2.208%
24	10.780	1576	1579	1582	rVB2	3930568	4739233	8.39%	0.695%
25	10.847	1582	1590	1594	rBV2	34655441	56476525	100.00%	8.286%
26	10.920	1594	1602	1606	rBV	34033912	56308450	99.70%	8.261%
27	10.963	1606	1609	1613	rVB	17803164	22190612	39.29%	3.256%
28	11.018	1613	1618	1623	rVB3	5643176	9885760	17.50%	1.450%
29	11.073	1623	1627	1628	rBV2	7356030	10183835	18.03%	1.494%
30	11.091	1628	1630	1636	rVB2	10059798	14100515	24.97%	2.069%
31	11.195	1643	1647	1650	rBV	6279429	7959345	14.09%	1.168%
32	11.225	1650	1652	1654	rVV	2740203	3039948	5.38%	0.446%
33	11.286	1654	1662	1671	rVB3	22602285	41382568	73.27%	6.071%
34	11.365	1671	1675	1678	rBV	8659786	11789920	20.88%	1.730%

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Integration Parameters: RTEINT.P

Integrator: RTE
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 Sampling : 1
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 Peak Location: TOP

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35	11.408	1678	1682	1686	rVV3	7447625	13282210	23.52%	1.949%
36	11.451	1686	1689	1693	rVV3	9709841	14473104	25.63%	2.123%
37	11.499	1693	1697	1701	rVV	20098123	28306293	50.12%	4.153%
38	11.542	1701	1704	1709	rVB3	13185275	19715975	34.91%	2.893%
39	11.591	1709	1712	1717	rVB2	3175865	4393363	7.78%	0.645%
40	11.640	1717	1720	1722	rBV	2700522	3217279	5.70%	0.472%
41	11.695	1722	1729	1733	rVB6	10016469	20081926	35.56%	2.946%
42	11.749	1733	1738	1740	rBV2	6622276	8692978	15.39%	1.275%
43	11.780	1740	1743	1746	rBV	23870627	27047754	47.89%	3.968%
44	11.810	1746	1748	1750	rVB2	1062919	790762	1.40%	0.116%
45	11.902	1759	1763	1765	rBV2	5985007	9155355	16.21%	1.343%
46	11.932	1765	1768	1769	rVV	6754727	8866658	15.70%	1.301%
47	11.957	1769	1772	1777	rVB2	12598854	18255756	32.32%	2.678%
48	12.012	1777	1781	1784	rBV	17335353	20592148	36.46%	3.021%
49	12.091	1791	1794	1799	rBV5	2879844	4790492	8.48%	0.703%
50	12.133	1799	1801	1804	rVB3	2002522	2252861	3.99%	0.331%
51	12.176	1804	1808	1810	rBV2	1648332	2367156	4.19%	0.347%
52	12.201	1810	1812	1814	rBV2	899515	893254	1.58%	0.131%
53	12.310	1825	1830	1835	rBV2	8458652	11389723	20.17%	1.671%
54	12.377	1836	1841	1848	rVV2	14793685	26312847	46.59%	3.861%
55	12.481	1852	1858	1863	rVV3	3830801	9268688	16.41%	1.360%
56	12.530	1863	1866	1871	rVB4	4008802	7068457	12.52%	1.037%
57	12.615	1874	1880	1882	rBV3	2935087	5195817	9.20%	0.762%
58	12.676	1886	1890	1893	rVB	5903595	7160912	12.68%	1.051%
59	12.768	1898	1905	1911	rVV5	3351211	9726569	17.22%	1.427%
60	12.822	1911	1914	1919	rVV4	1265801	1990633	3.52%	0.292%
61	12.883	1919	1924	1930	rVB4	4944438	8476871	15.01%	1.244%
62	12.950	1930	1935	1938	rBV3	3546020	5821327	10.31%	0.854%
63	12.981	1938	1940	1945	rVB2	3842361	4270164	7.56%	0.627%
64	13.048	1947	1951	1954	rVV2	2814880	3424557	6.06%	0.502%
65	13.085	1954	1957	1964	rVB2	1438073	2571485	4.55%	0.377%
66	13.152	1964	1968	1973	rBV2	1036825	1817724	3.22%	0.267%
67	13.200	1973	1976	1979	rVB	698565	725569	1.28%	0.106%
68	13.347	1996	2000	2007	rVB2	3578466	5501725	9.74%	0.807%
69	13.475	2017	2021	2025	rVB6	592249	892680	1.58%	0.131%
70	13.639	2045	2048	2052	rVV2	533213	687169	1.22%	0.101%
71	13.694	2052	2057	2060	rVV4	281391	636594	1.13%	0.093%

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ClientSampleId :
EP1-MW-A(8-12)

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 >
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72	13.731	2060	2063	2071	rVB6	383594	726738	1.29%	0.107%
73	13.810	2071	2076	2081	rBV3	321644	580655	1.03%	0.085%

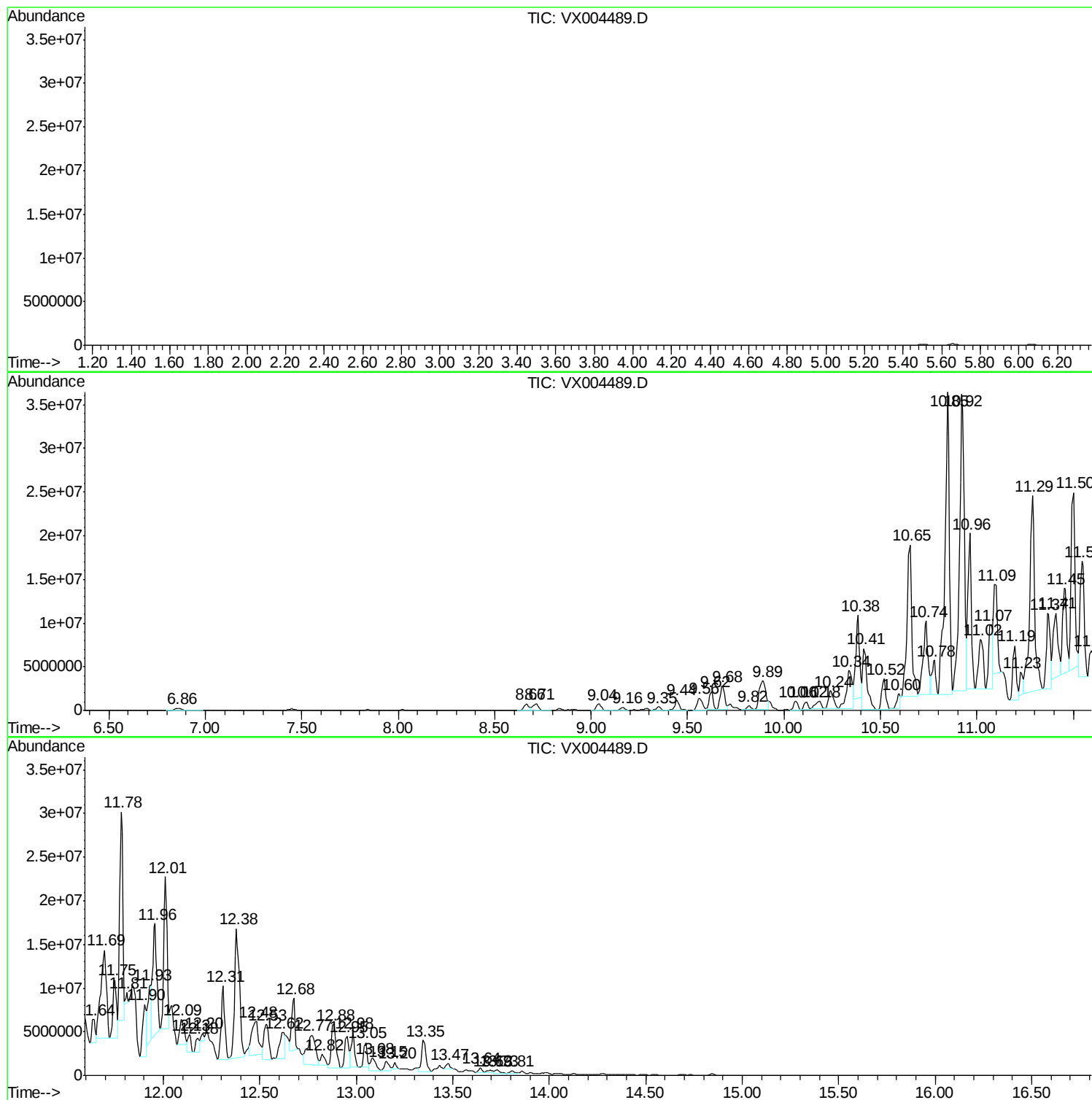
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Instrument :
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ClientSampleId :
EP1-MW-A(8-12)

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



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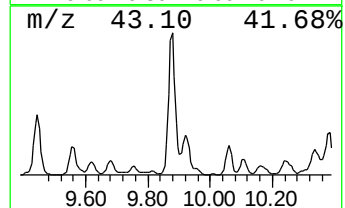
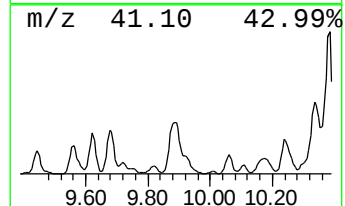
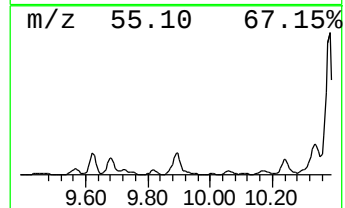
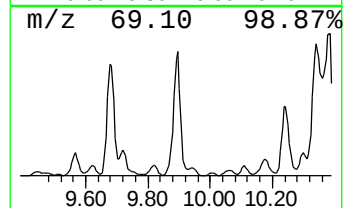
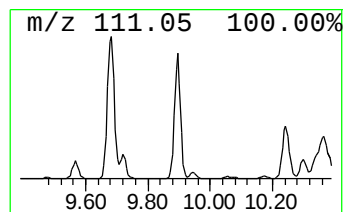
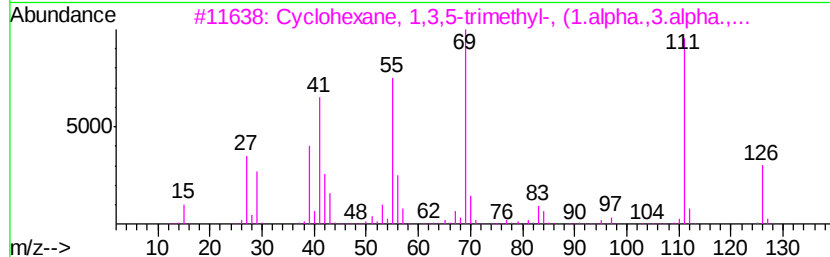
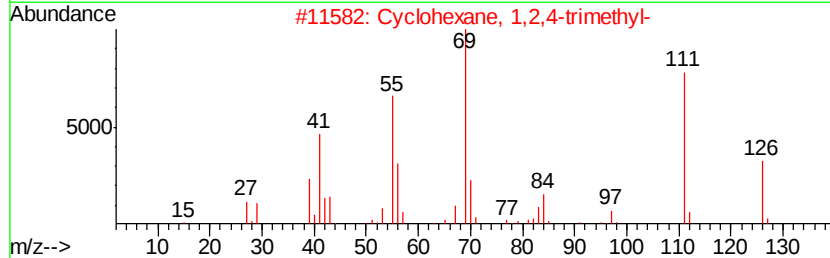
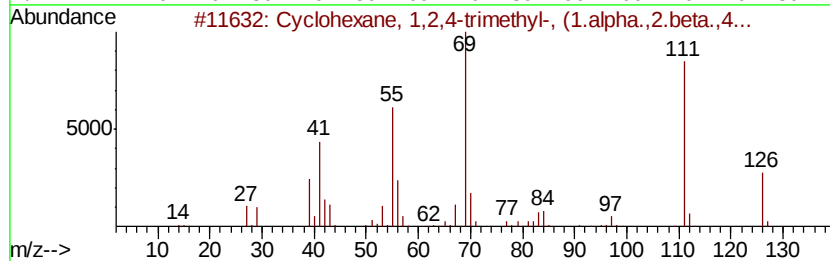
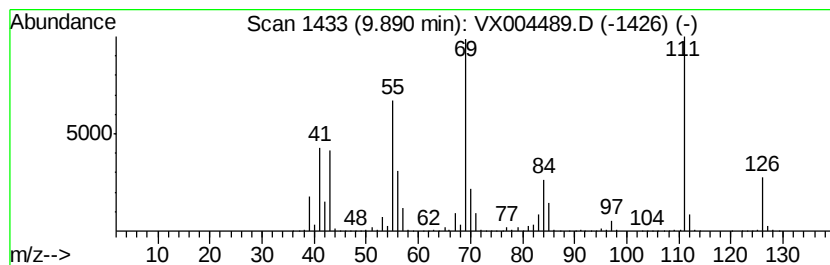
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, 1,2,4-trimethy... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.89	319.62 ug/l	8071360	Chlorobenzene-d5	10.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,2,4-trimethyl-, (...)	126	C9H18	007667-60-9	91
2		Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	83
3		Cyclohexane, 1,3,5-trimethyl-, (...)	126	C9H18	001795-26-2	74
4		Cyclohexane, 1,3,5-trimethyl-, (...)	126	C9H18	001795-27-3	68
5		Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	68



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Instrument :
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ClientSampled :
EP1-MW-A(8-12)

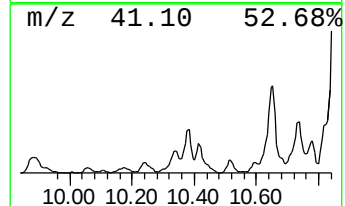
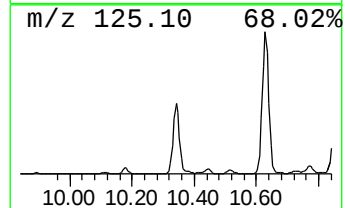
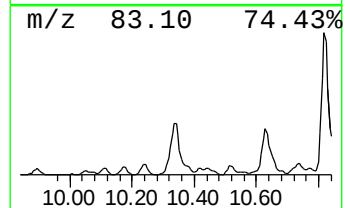
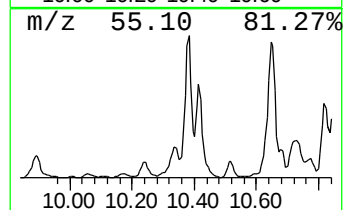
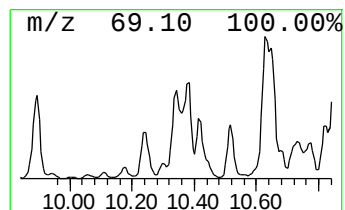
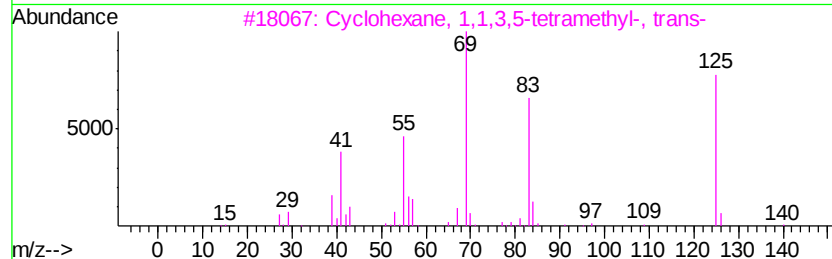
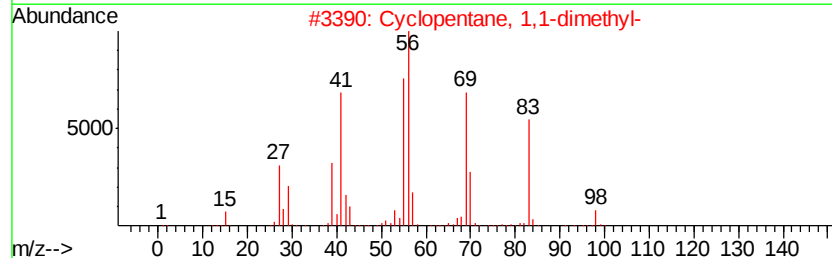
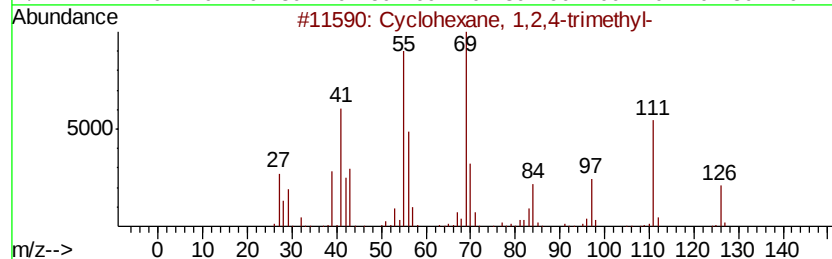
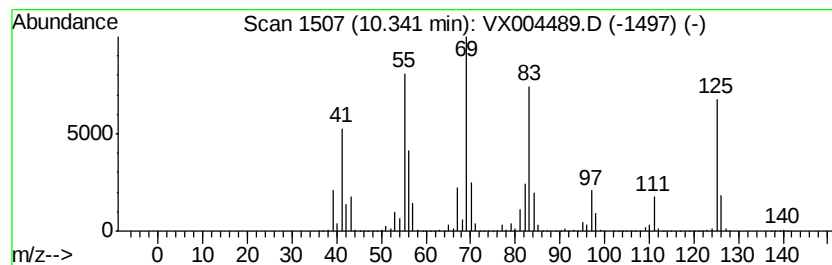
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexane, 1,2,4-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.34	371.42 ug/l	9379490	Chlorobenzene-d5	10.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	70
2		Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	45
3		Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-31-8	43
4		3-Heptene, 2,6-dimethyl-	126	C9H18	002738-18-3	41
5		cis-3-Nonene	126	C9H18	020237-46-1	38



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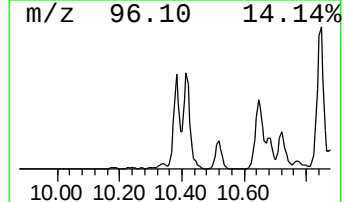
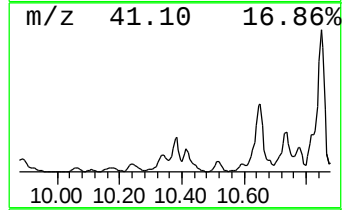
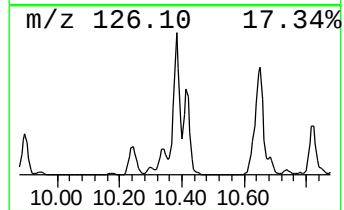
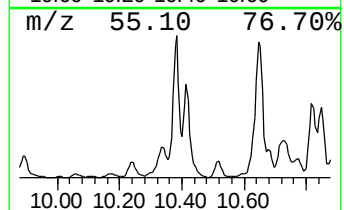
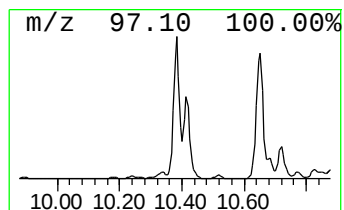
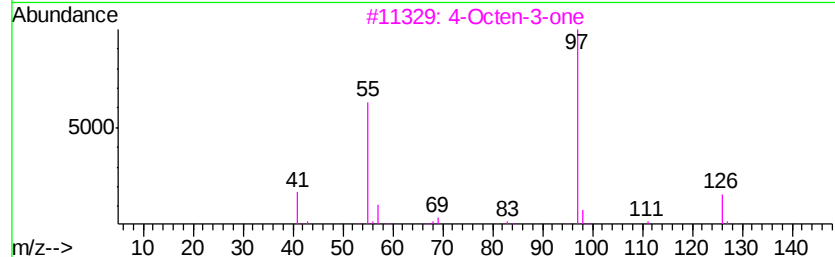
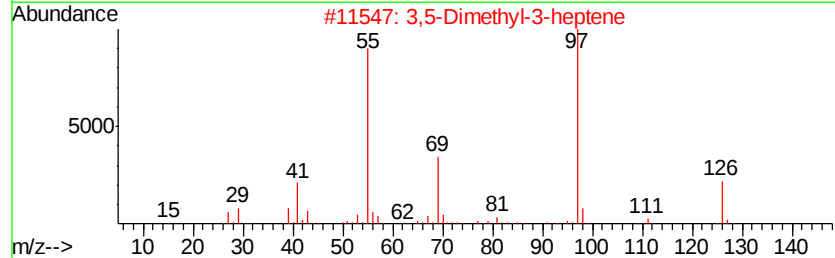
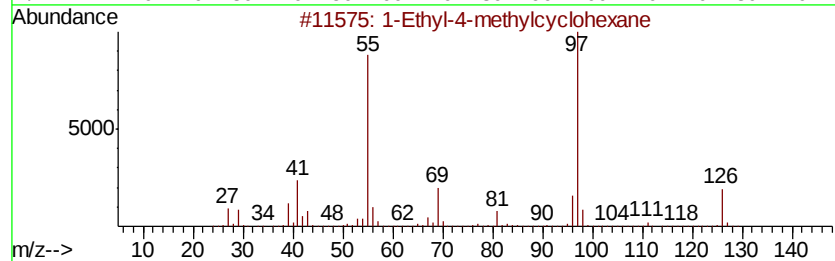
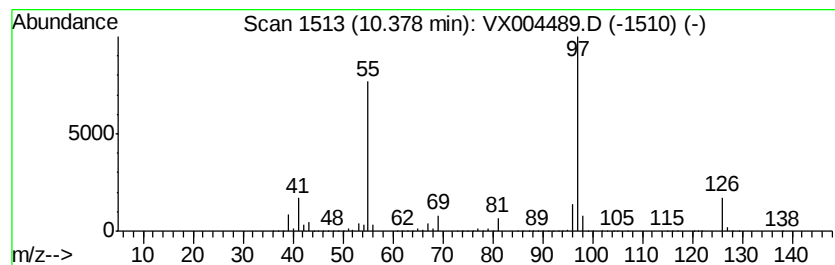
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Quant Title : SW846 8260

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

Peak Number 3 1-Ethyl-4-methylcyclohexane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.38	519.83 ug/l	13127400	Chlorobenzene-d5	10.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	91
2		3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	80
3		4-Octen-3-one	126	C8H14O	014129-48-7	80
4		Cyclohexane, 1-bromo-4-methyl-	176	C7H13Br	006294-40-2	64
5		Cycloheptane, bromo-	176	C7H13Br	002404-35-5	59



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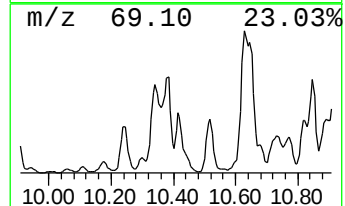
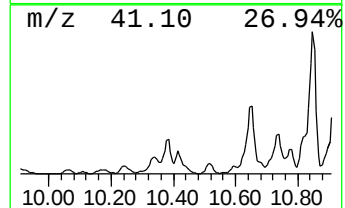
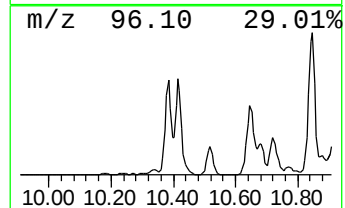
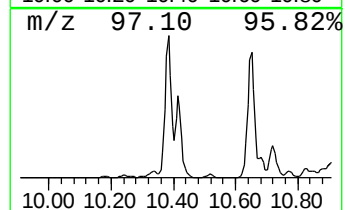
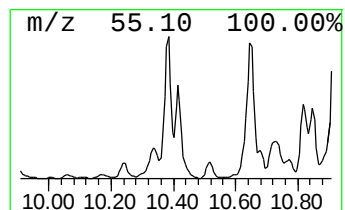
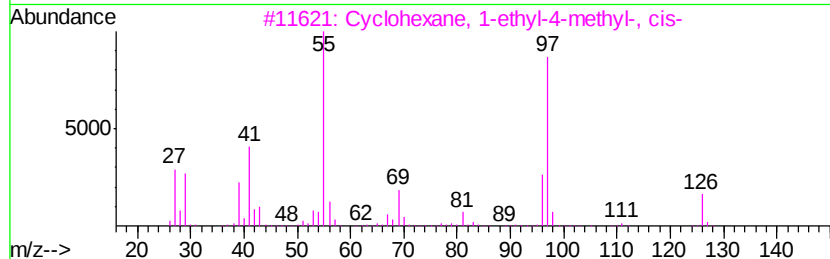
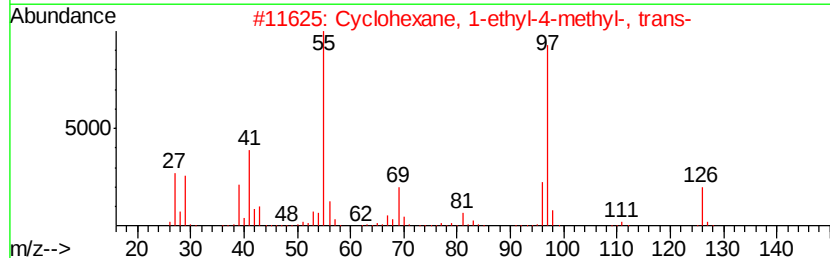
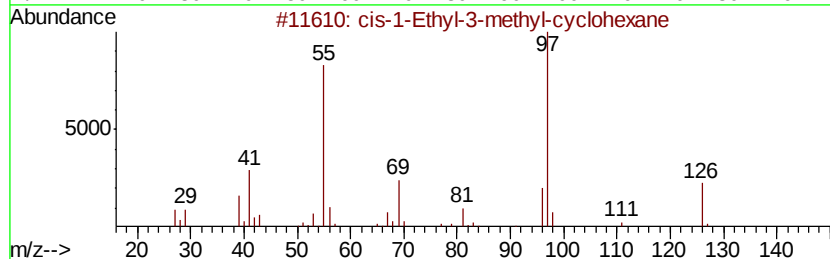
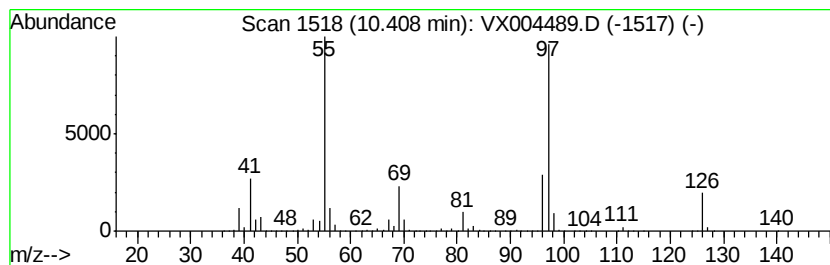
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Quant Title : SW846 8260

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

Peak Number 4 cis-1-Ethyl-3-methyl-cyclohexane... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.41	438.97 ug/l	11085300	Chlorobenzene-d5	10.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	93
2		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	91
3		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	91
4		1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	90
5		1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091218\
Data File : VX004489.D
Acq On : 12 Sep 2018 12:46
Operator : JC/MD
Sample : J4848-02
Misc : 2.00µL/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EP1-MW-A(8-12)

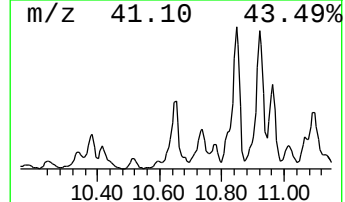
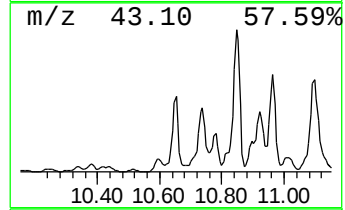
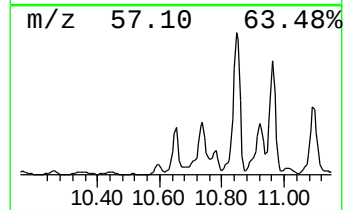
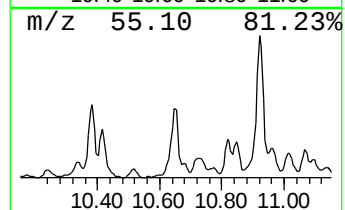
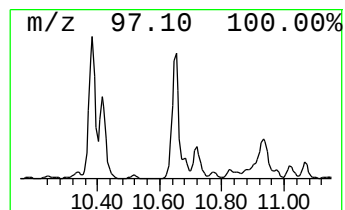
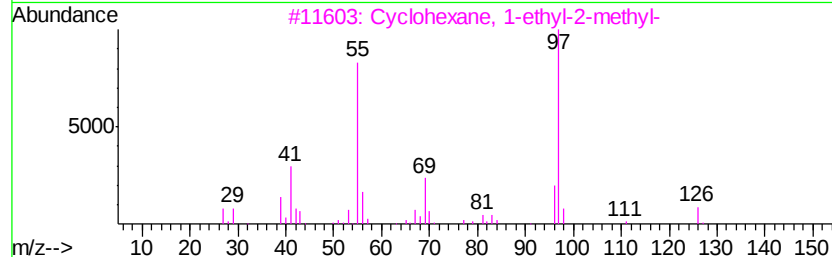
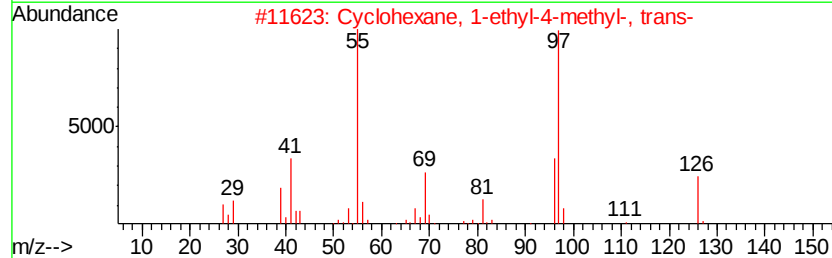
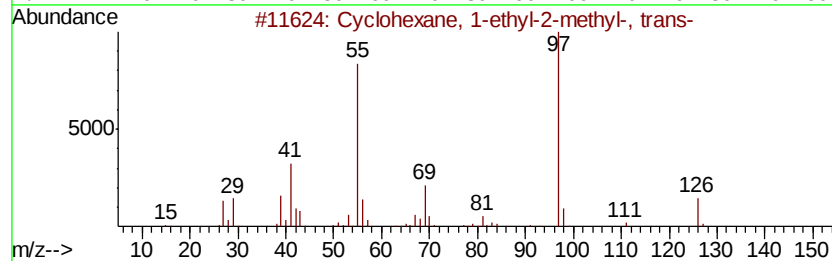
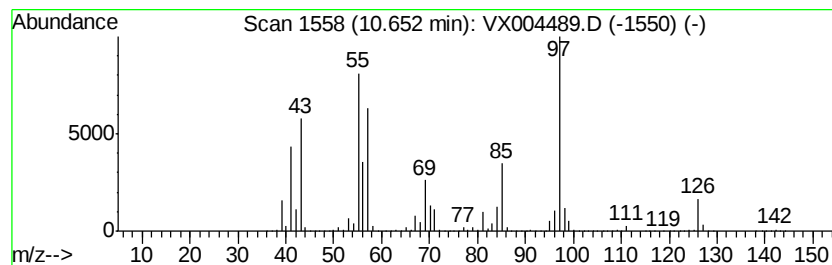
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Quant Title : SW846 8260

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

Peak Number 5 Cyclohexane, 1-ethyl-2-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	1204.38 ug/l	30414500	Chlorobenzene-d5	10.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	70
2		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	64
3		Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	62
4		1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	58
5		3-Heptene, 3-ethyl-	126	C9H18	074764-46-8	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091218\
Data File : VX004489.D
Acq On : 12 Sep 2018 12:46
Operator : JC/MD
Sample : J4848-02
Misc : 2.00µL/5mL/100µL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

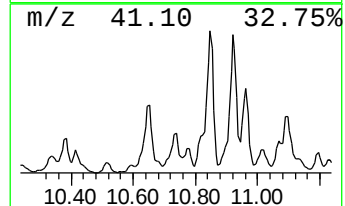
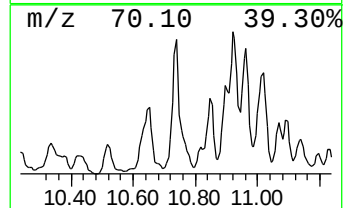
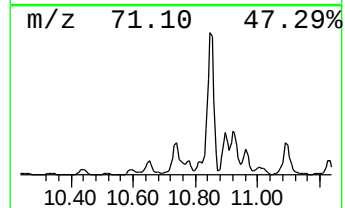
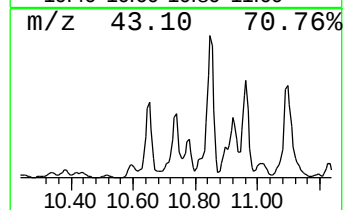
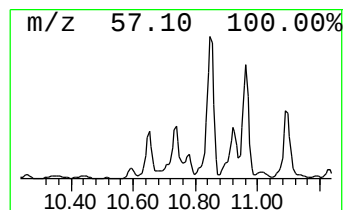
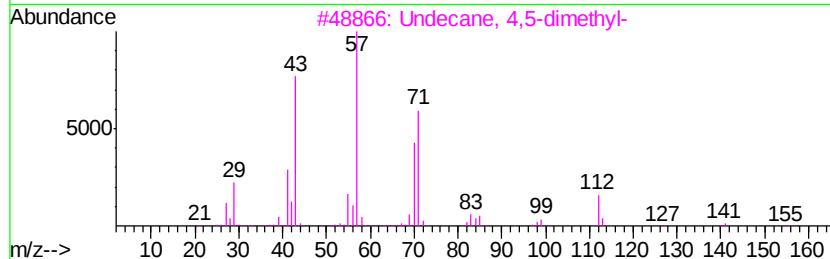
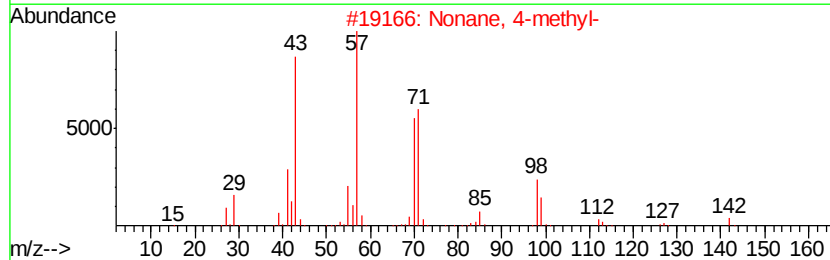
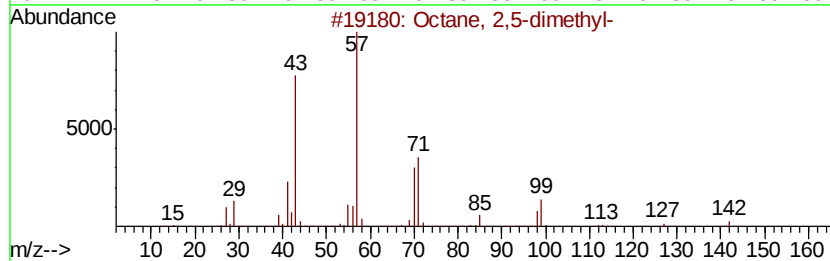
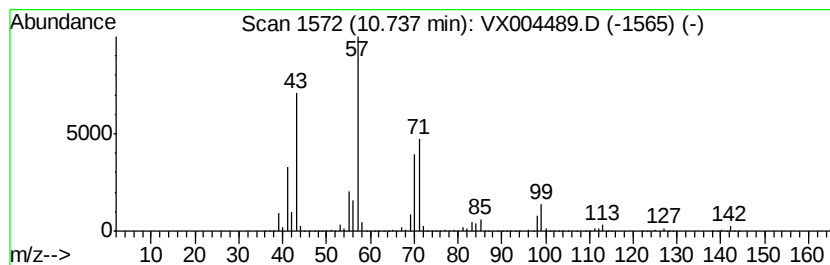
Instrument :
MSVOA_X
ClientSampled :
EP1-MW-A(8-12)

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 Octane, 2,5-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.74	596.01 ug/l	15051200	Chlorobenzene-d5	10.12	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,5-dimethyl-	142	C10H22	015869-89-3	91
2	Nonane, 4-methyl-	142	C10H22	017301-94-9	80
3	Undecane, 4,5-dimethyl-	184	C13H28	017312-79-7	78
4	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	58
5	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091218\
Data File : VX004489.D
Acq On : 12 Sep 2018 12:46
Operator : JC/MD
Sample : J4848-02
Misc : 2.00uL/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

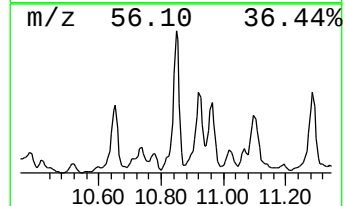
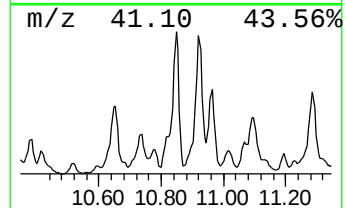
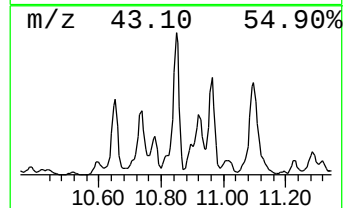
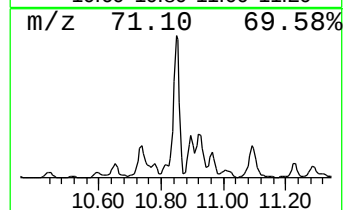
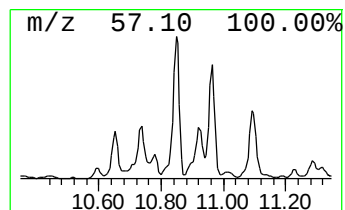
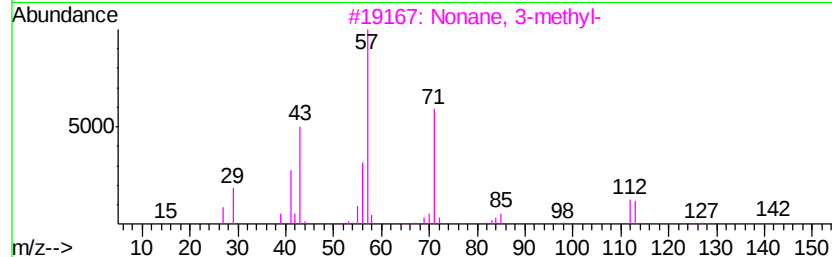
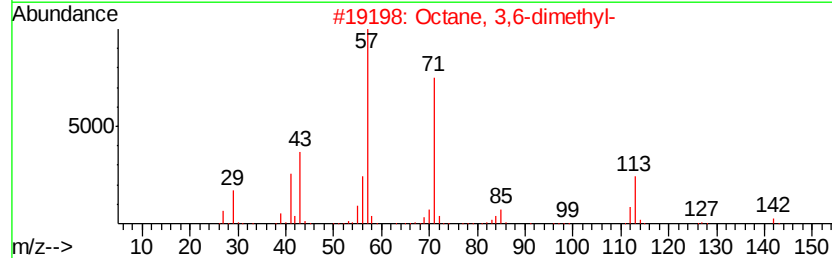
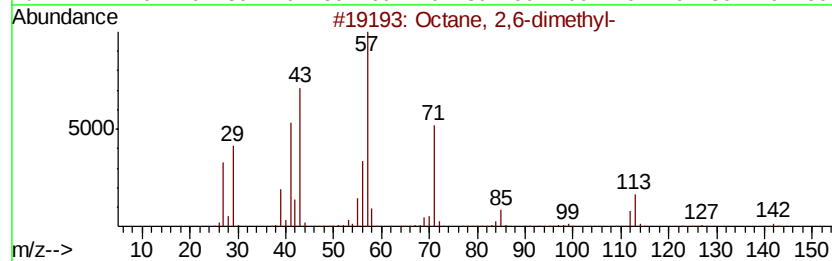
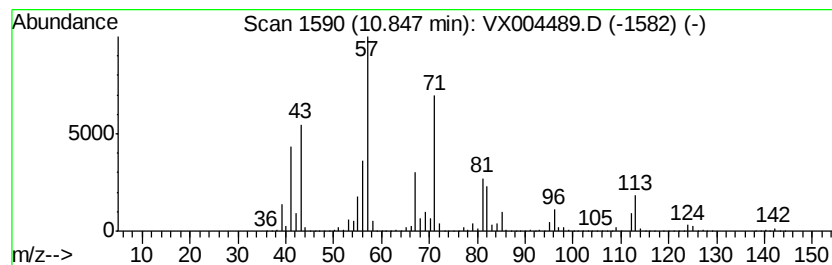
Instrument :
MSVOA_X
ClientSampleId :
EP1-MW-A(8-12)

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 Octane, 2,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.85	2236.41 ug/l	56476500	Chlorobenzene-d5	10.12
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Octane, 2,6-dimethyl-	142 C10H22	002051-30-1 70
2		Octane, 3,6-dimethyl-	142 C10H22	015869-94-0 62
3		Nonane, 3-methyl-	142 C10H22	005911-04-6 60
4		Sulfurous acid, 2-ethylhexyl hex...	418 C24H50O3S	1000309-19-9 47
5		Undecane, 6,6-dimethyl-	184 C13H28	017312-76-4 43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091218\
Data File : VX004489.D
Acq On : 12 Sep 2018 12:46
Operator : JC/MD
Sample : J4848-02
Misc : 2.00µL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
EP1-MW-A(8-12)

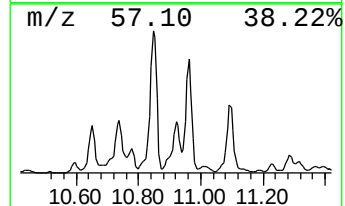
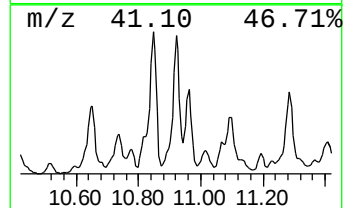
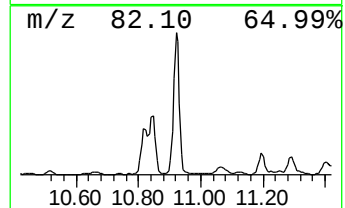
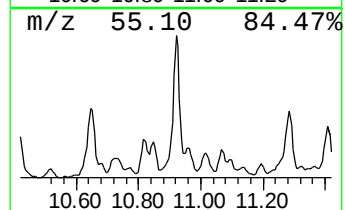
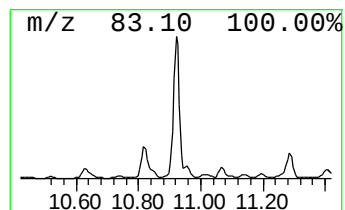
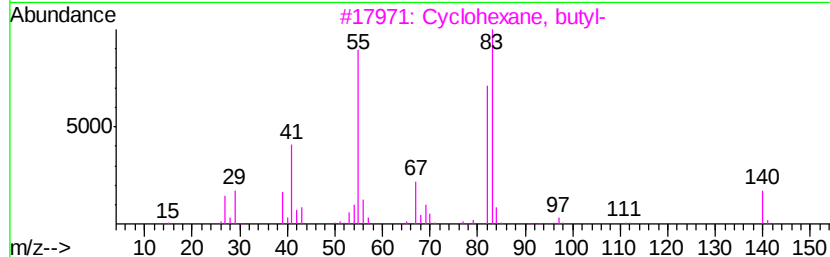
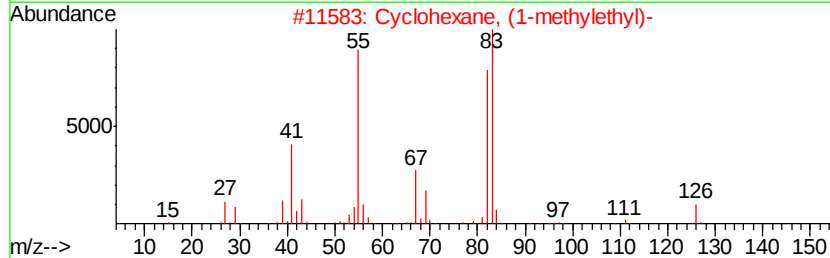
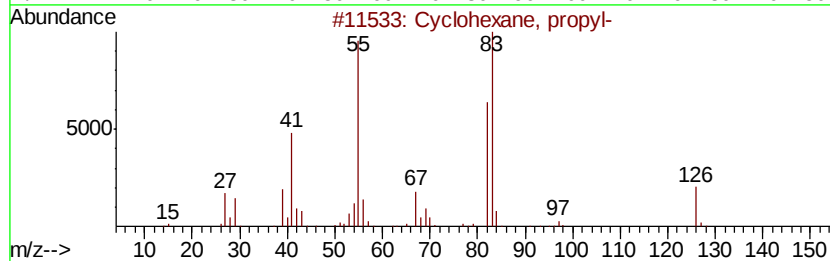
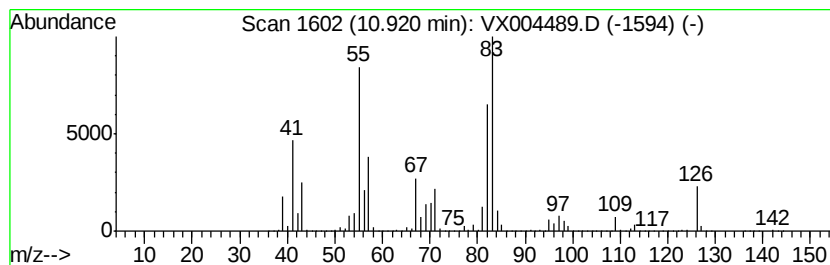
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Quant Title : SW846 8260

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

Peak Number 8 Cyclohexane, propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.92	2229.75 ug/l	56308500	Chlorobenzene-d5	10.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, propyl-	126	C9H18	001678-92-8	81
2		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	64
3		Cyclohexane, butyl-	140	C10H20	001678-93-9	59
4		Cyclohexane, octyl-	196	C14H28	001795-15-9	59
5		Cyclohexane, ethyl-	112	C8H16	001678-91-7	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091218\
Data File : VX004489.D
Acq On : 12 Sep 2018 12:46
Operator : JC/MD
Sample : J4848-02
Misc : 2.00µL/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

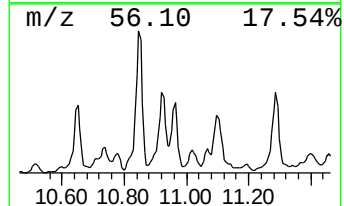
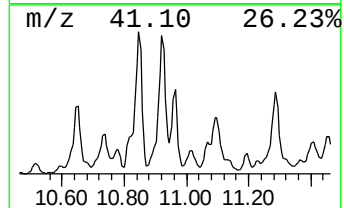
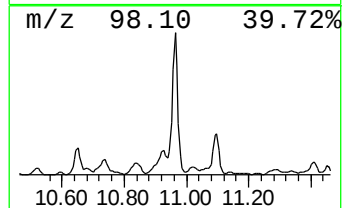
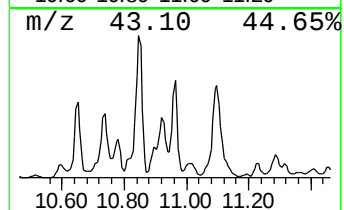
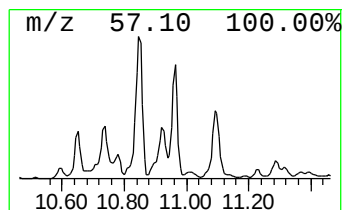
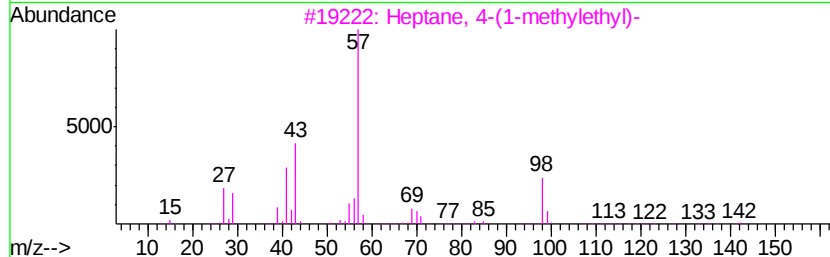
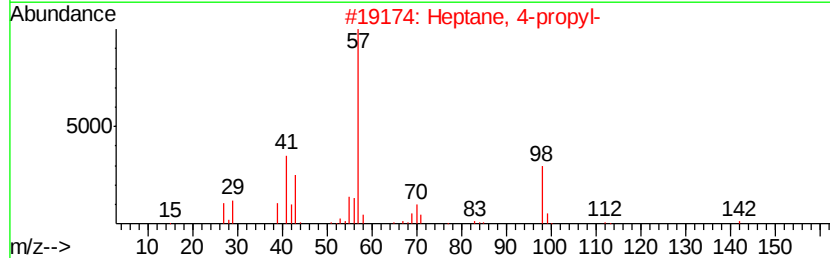
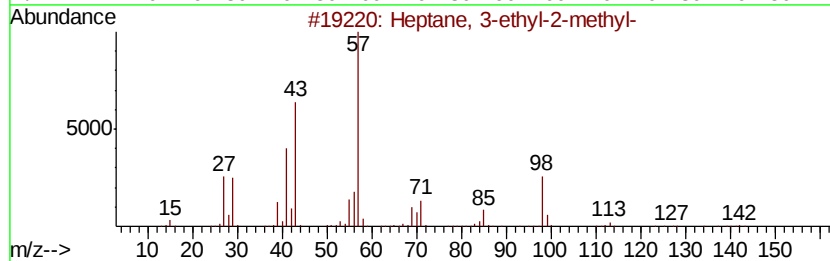
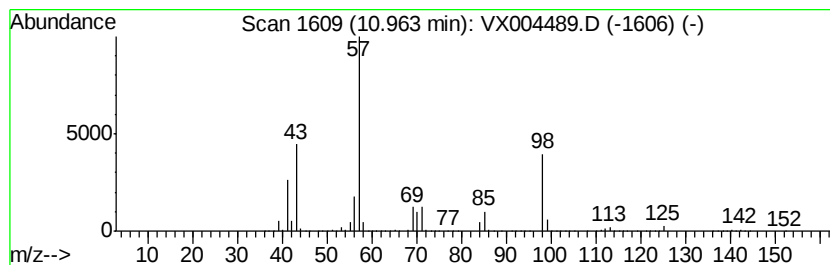
Instrument :
MSVOA_X
ClientSampled :
EP1-MW-A(8-12)

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 9 Heptane, 3-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.96	878.72 ug/l	22190600	Chlorobenzene-d5	10.12
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptane, 3-ethyl-2-methyl-	142 C10H22	014676-29-0	74
2	Heptane, 4-propyl-	142 C10H22	003178-29-8	64
3	Heptane, 4-(1-methylethyl)-	142 C10H22	052896-87-4	52
4	Hexane, 3-ethyl-2,5-dimethyl-	142 C10H22	052897-04-8	50
5	Undecane, 6-methyl-	170 C12H26	017302-33-9	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091218\
Data File : VX004489.D
Acq On : 12 Sep 2018 12:46
Operator : JC/MD
Sample : J4848-02
Misc : 2.00µL/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

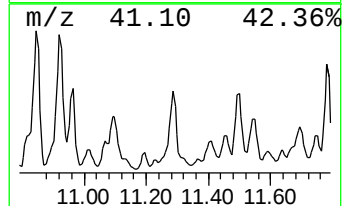
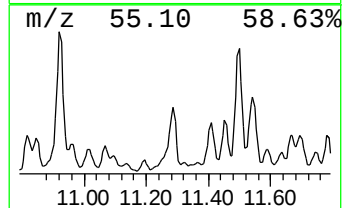
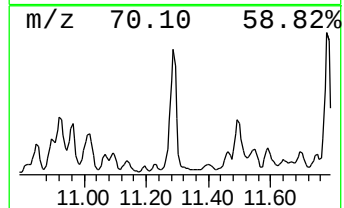
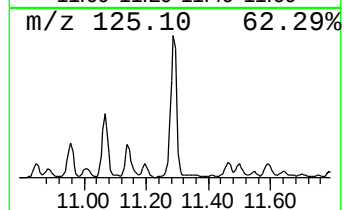
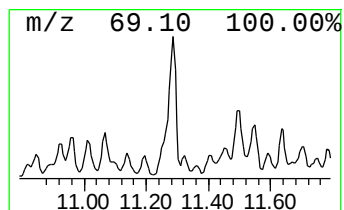
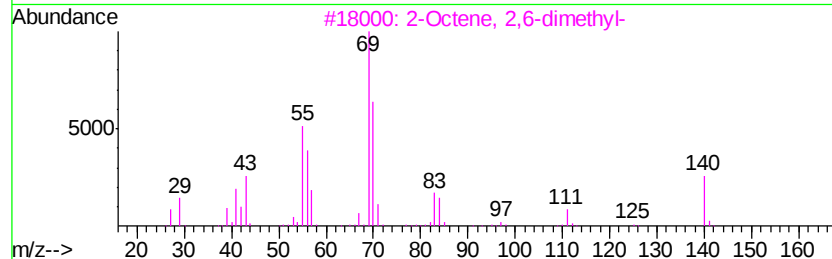
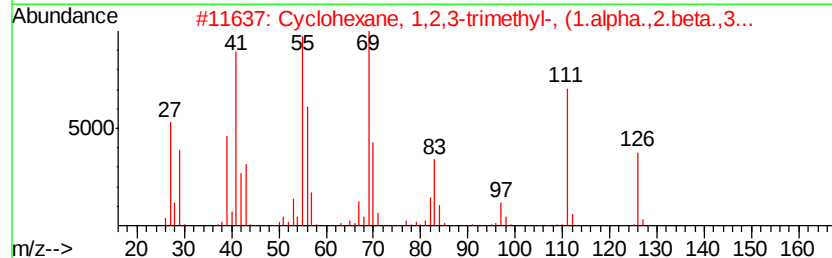
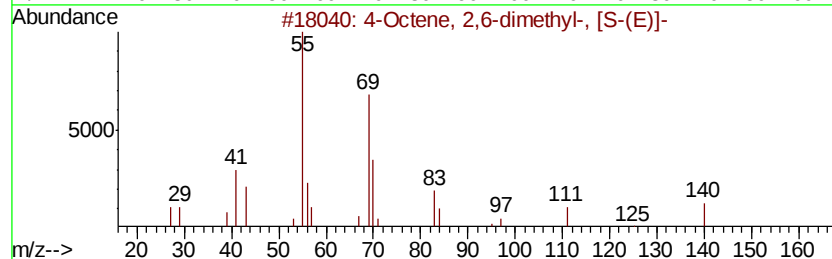
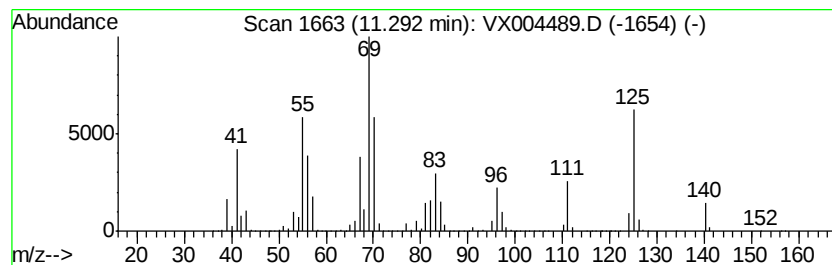
Instrument :
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ClientSampleId :
EP1-MW-A(8-12)

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 4-Octene, 2,6-dimethyl-, [S... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.29	431.92 ug/l	41382600	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	4-Octene, 2,6-dimethyl-, [S-(E)]-	140 C10H20	062960-76-3	62
2	Cyclohexane, 1,2,3-trimethyl-, (...)	126 C9H18	001678-81-5	60
3	2-Octene, 2,6-dimethyl-	140 C10H20	004057-42-5	58
4	4-Octene, 2,6-dimethyl-, [S-(Z)]-	140 C10H20	062960-77-4	58
5	Cyclohexane, 1,1,2,3-tetramethyl-	140 C10H20	006783-92-2	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX091218\
Data File : VX004489.D
Acq On : 12 Sep 2018 12:46
Operator : JC/MD
Sample : J4848-02
Misc : 2.00g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EP1-MW-A(8-12)

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclohexane, 1,2,...	9.89	319.6	ug/l	8071360	3	10.12	1262660	50.0
Cyclohexane, 1,2,...	10.34	371.4	ug/l	9379490	3	10.12	1262660	50.0
1-Ethyl-4-methylc...	10.38	519.8	ug/l	13127400	3	10.12	1262660	50.0
cis-1-Ethyl-3-met...	10.41	439.0	ug/l	11085300	3	10.12	1262660	50.0
Cyclohexane, 1-et...	10.65	1204.4	ug/l	30414500	3	10.12	1262660	50.0
Octane, 2,5-dimet...	10.74	596.0	ug/l	15051200	3	10.12	1262660	50.0
Octane, 2,6-dimet...	10.85	2236.4	ug/l	56476500	3	10.12	1262660	50.0
Cyclohexane, propyl-	10.92	2229.8	ug/l	56308500	3	10.12	1262660	50.0
Heptane, 3-ethyl-...	10.96	878.7	ug/l	22190600	3	10.12	1262660	50.0
4-Octene, 2,6-dim...	11.29	431.9	ug/l	41382600	4	12.09	4790490	50.0