

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
 Data File : VX006009.D
 Acq On : 15 Nov 2018 07:39
 Operator : JC/MD
 Sample : J5915-07
 Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 T-7

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\82X110718W.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	2.843	272	277	288	rVB	50739	104272	9.09%	0.372%
2	3.129	315	324	335	rBV	30880	70201	6.12%	0.250%
3	3.483	373	382	396	rVV2	26890	62880	5.48%	0.224%
4	4.367	519	527	540	rVB2	21341	62558	5.45%	0.223%
5	5.507	694	714	721	rBV2	68341	226333	19.72%	0.807%
6	5.586	721	727	732	rVV2	51424	142878	12.45%	0.510%
7	5.659	732	739	765	rVB2	138910	507359	44.21%	1.810%
8	5.903	768	779	794	rVB	66527	190980	16.64%	0.681%
9	6.068	795	806	823	rBV2	78088	220112	19.18%	0.785%
10	6.336	840	850	860	rVB3	103740	303581	26.45%	1.083%
11	6.689	896	908	916	rBV2	72745	177315	15.45%	0.632%
12	6.860	927	936	943	rBV2	192702	488270	42.55%	1.742%
13	7.445	1021	1032	1044	rVB4	49145	136609	11.90%	0.487%
14	7.567	1044	1052	1057	rBV	41402	87454	7.62%	0.312%
15	7.628	1057	1062	1073	rVV2	53090	117607	10.25%	0.419%
16	8.049	1121	1131	1143	rVB	95465	209028	18.21%	0.746%
17	8.171	1143	1151	1159	rBV2	84115	211944	18.47%	0.756%
18	8.250	1159	1164	1173	rVV2	69275	151125	13.17%	0.539%
19	8.336	1173	1178	1182	rVV	161748	284015	24.75%	1.013%
20	8.372	1182	1184	1191	rVV	80253	116271	10.13%	0.415%
21	8.494	1199	1204	1213	rVV	180006	353980	30.85%	1.263%
22	8.671	1226	1233	1236	rBV2	89126	181567	15.82%	0.648%
23	8.713	1236	1240	1249	rVV	375217	621544	54.16%	2.217%
24	8.829	1249	1259	1264	rVV4	31316	80448	7.01%	0.287%
25	8.878	1264	1267	1276	rVV2	32279	80994	7.06%	0.289%
26	8.970	1276	1282	1290	rVV	210549	354078	30.85%	1.263%
27	9.043	1290	1294	1303	rVV4	29443	77593	6.76%	0.277%
28	9.256	1326	1329	1339	rVB4	27611	62946	5.48%	0.225%
29	9.439	1354	1359	1365	rBV	62457	100133	8.73%	0.357%
30	9.555	1373	1378	1385	rBV2	113900	222297	19.37%	0.793%
31	9.817	1415	1421	1426	rBV2	35003	68144	5.94%	0.243%
32	9.872	1426	1430	1434	rVV2	53351	97364	8.48%	0.347%
33	9.921	1435	1438	1440	rVV2	62079	89544	7.80%	0.319%
34	9.957	1440	1444	1449	rVB	390939	539049	46.97%	1.923%

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35	10.061	1456	1461	1466	rBV	310444	415661	36.22%	1.483%
36	10.116	1466	1470	1474	rBV	368913	508165	44.28%	1.812%
37	10.256	1487	1493	1497	rVB	223906	321801	28.04%	1.148%
38	10.317	1497	1503	1504	rVV4	35406	62542	5.45%	0.223%
39	10.360	1505	1510	1515	rVV	281956	444286	38.71%	1.585%
40	10.414	1515	1519	1530	rVB	346879	549520	47.88%	1.960%
41	10.640	1552	1556	1560	rBV	119835	149663	13.04%	0.534%
42	10.725	1566	1570	1575	rBV	80039	111544	9.72%	0.398%
43	10.835	1580	1588	1592	rBV	126951	214087	18.66%	0.764%
44	10.914	1592	1601	1605	rBV3	87741	174818	15.23%	0.624%
45	11.018	1612	1618	1623	rBV	90164	154615	13.47%	0.551%
46	11.079	1624	1628	1632	rBV	87641	131169	11.43%	0.468%
47	11.140	1632	1638	1641	rBV	454002	755000	65.79%	2.693%
48	11.244	1652	1655	1659	rVV	231126	263601	22.97%	0.940%
49	11.359	1670	1674	1682	rVV	334291	458774	39.98%	1.636%
50	11.439	1682	1687	1689	rVV	218515	322687	28.12%	1.151%
51	11.506	1693	1698	1699	rVV2	143441	205941	17.95%	0.735%
52	11.530	1699	1702	1710	rVB	362381	478169	41.67%	1.705%
53	11.670	1721	1725	1732	rBV2	172991	279878	24.39%	0.998%
54	11.768	1738	1741	1744	rVV	62380	60559	5.28%	0.216%
55	11.810	1744	1748	1756	rVB2	459076	670486	58.42%	2.391%
56	11.890	1757	1761	1764	rBV	70672	104301	9.09%	0.372%
57	11.945	1768	1770	1774	rVB2	87856	98855	8.61%	0.353%
58	11.987	1774	1777	1781	rBV2	73347	109729	9.56%	0.391%
59	12.079	1786	1792	1797	rBV2	458579	685921	59.77%	2.446%
60	12.152	1797	1804	1807	rVV3	207974	434202	37.84%	1.549%
61	12.201	1810	1812	1814	rVV	58772	72819	6.35%	0.260%
62	12.225	1814	1816	1823	rVB2	144532	172859	15.06%	0.617%
63	12.304	1823	1829	1836	rBV	655295	1072381	93.44%	3.825%
64	12.371	1836	1840	1847	rVB2	438183	806431	70.27%	2.876%
65	12.475	1847	1857	1861	rBV2	292234	460981	40.17%	1.644%
66	12.524	1861	1865	1871	rVB2	206754	315510	27.49%	1.125%
67	12.591	1871	1876	1878	rBV	171978	243641	21.23%	0.869%
68	12.609	1878	1879	1882	rVV	110849	103527	9.02%	0.369%
69	12.670	1885	1889	1893	rVV	979288	1147607	100.00%	4.093%
70	12.701	1893	1894	1898	rVB	80792	67587	5.89%	0.241%
71	12.755	1898	1903	1912	rVB3	244057	577239	50.30%	2.059%

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72	12.853	1913	1919	1924	rBV5	42151	93332	8.13%	0.333%
73	12.902	1924	1927	1931	rVB2	62168	77007	6.71%	0.275%
74	12.981	1931	1940	1944	rBV	519212	833487	72.63%	2.973%
75	13.024	1944	1947	1952	rVB2	223392	291668	25.42%	1.040%
76	13.079	1952	1956	1958	rBV	186372	257824	22.47%	0.920%
77	13.127	1962	1964	1967	rVB	97545	90373	7.87%	0.322%
78	13.225	1973	1980	1984	rVB2	363022	541771	47.21%	1.932%
79	13.268	1984	1987	1990	rVB2	108785	120705	10.52%	0.431%
80	13.316	1990	1995	1997	rBV3	283120	391308	34.10%	1.396%
81	13.353	1997	2001	2006	rVV3	664459	1021492	89.01%	3.643%
82	13.426	2010	2013	2018	rVB	212563	262756	22.90%	0.937%
83	13.481	2018	2022	2030	rVB3	119975	216373	18.85%	0.772%
84	13.560	2030	2035	2038	rBV2	104289	156631	13.65%	0.559%
85	13.603	2038	2042	2044	rBV	166883	194081	16.91%	0.692%
86	13.633	2044	2047	2054	rVB2	269146	423595	36.91%	1.511%
87	13.731	2058	2063	2068	rVB3	173124	300276	26.17%	1.071%
88	13.835	2075	2080	2086	rVB	210249	289365	25.21%	1.032%
89	13.987	2098	2105	2109	rBV3	129265	226250	19.71%	0.807%
90	14.030	2109	2112	2115	rVV	82150	103241	9.00%	0.368%
91	14.091	2119	2122	2125	rVB3	125861	138071	12.03%	0.492%
92	14.133	2125	2129	2133	rBV	206575	232903	20.29%	0.831%
93	14.273	2148	2152	2162	rVB	186623	308979	26.92%	1.102%
94	14.377	2165	2169	2174	rBV	101565	126978	11.06%	0.453%
95	14.432	2174	2178	2182	rVB	111112	130795	11.40%	0.467%
96	14.542	2191	2196	2201	rBV5	36445	87057	7.59%	0.311%
97	14.694	2213	2221	2226	rBV	483058	629477	54.85%	2.245%
98	14.840	2242	2245	2249	rVB	213986	247076	21.53%	0.881%
99	15.554	2356	2362	2372	rVB2	79252	138065	12.03%	0.492%
100	15.688	2379	2384	2388	rBV	59930	99080	8.63%	0.353%

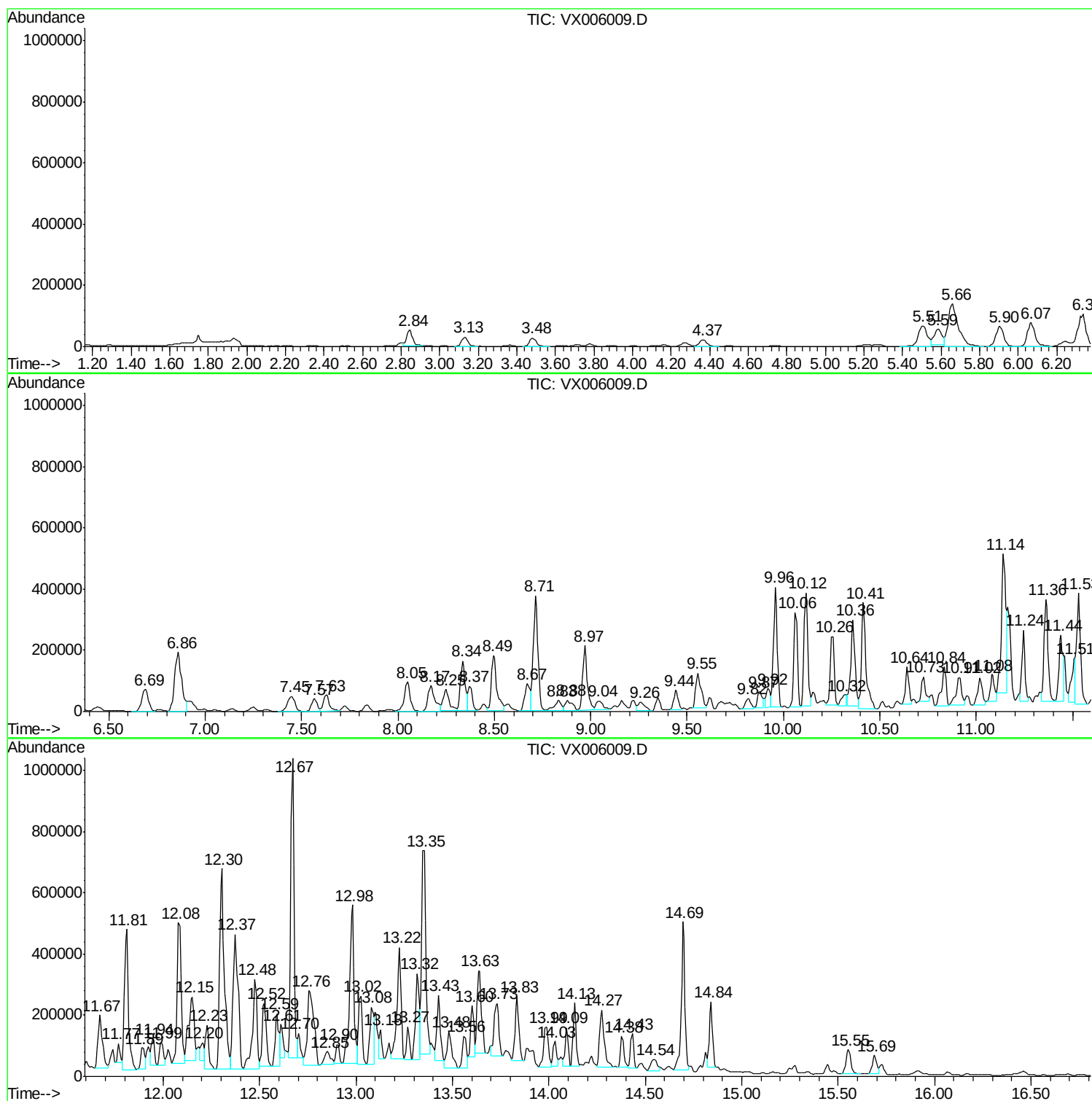
Sum of corrected areas: 28037045

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



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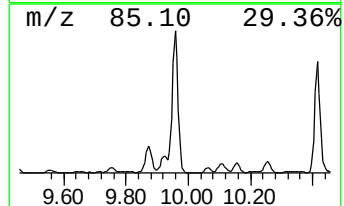
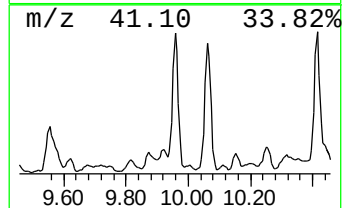
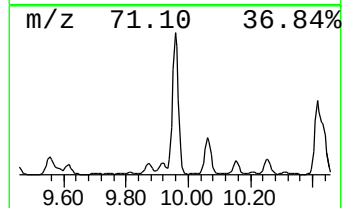
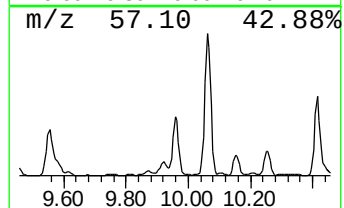
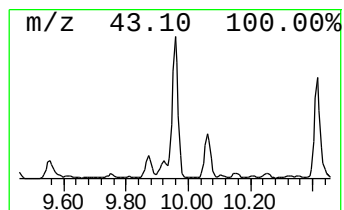
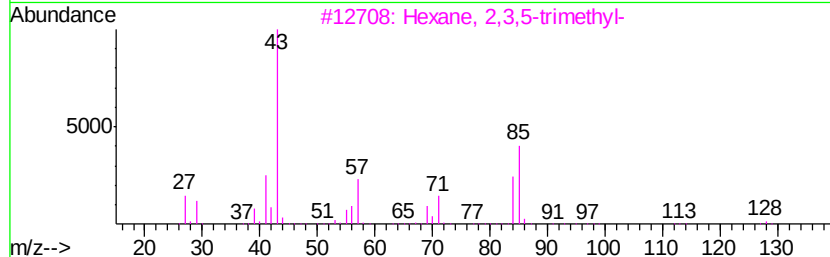
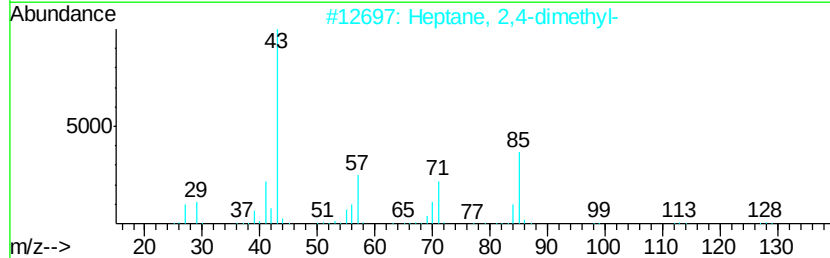
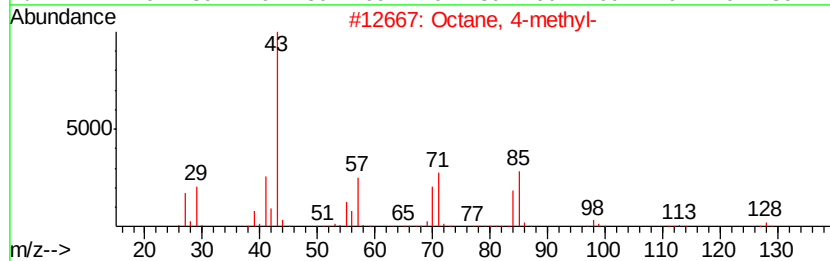
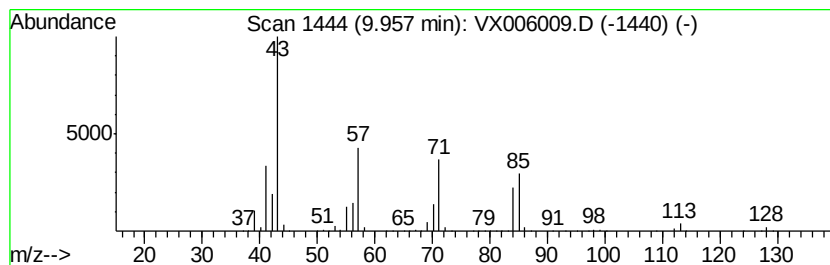
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

Peak Number 1 Octane, 4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.96	53.04 ug/l	539049	Chlorobenzene-d5	10.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 4-methyl-	128	C9H20	002216-34-4	80
2		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	64
3		Hexane, 2,3,5-trimethyl-	128	C9H20	001069-53-0	64
4		Ether, hexyl pentyl	172	C11H24O	032357-83-8	59
5		Decane, 2,4-dimethyl-	170	C12H26	002801-84-5	59



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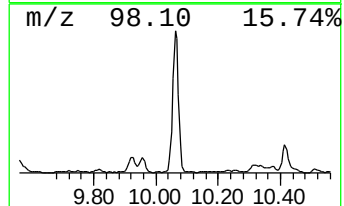
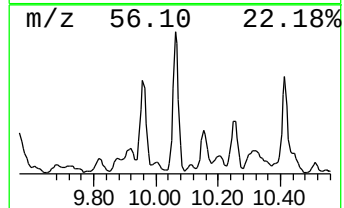
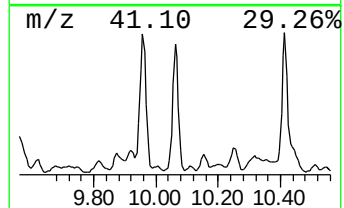
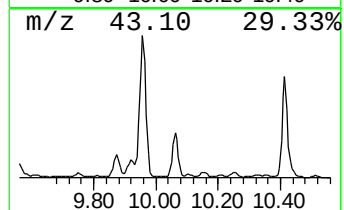
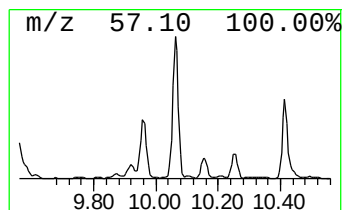
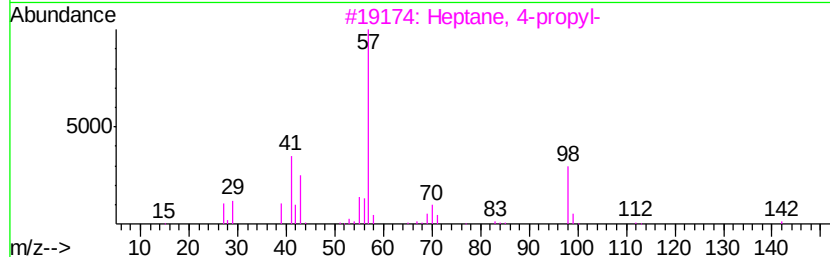
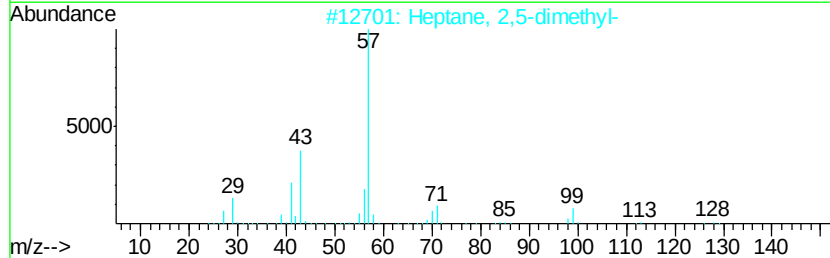
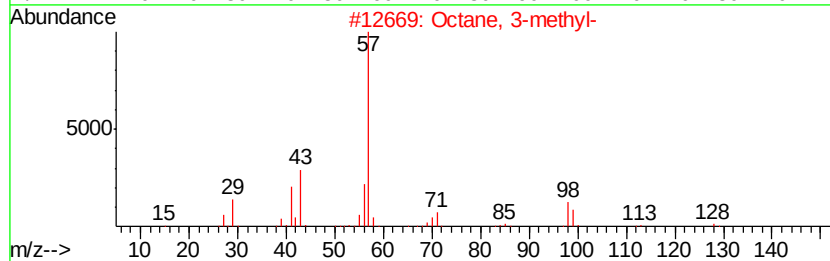
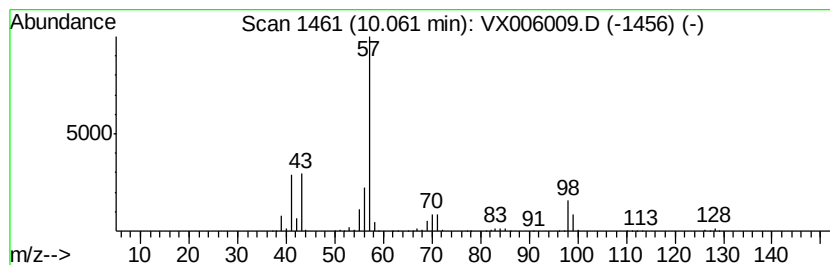
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Peak Number 2 Octane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.06	40.90 ug/l	415661	Chlorobenzene-d5	10.12
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Octane, 3-methyl-	128 C9H20	002216-33-3 90
2		Heptane, 2,5-dimethyl-	128 C9H20	002216-30-0 72
3		Heptane, 4-propyl-	142 C10H22	003178-29-8 64
4		Undecane, 6-methyl-	170 C12H26	017302-33-9 59
5		Oxalic acid, isobutyl heptyl ester	244 C13H24O4	1000309-37-2 59



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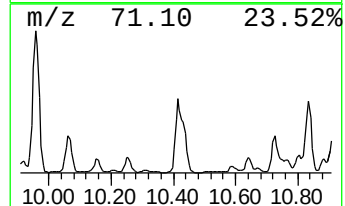
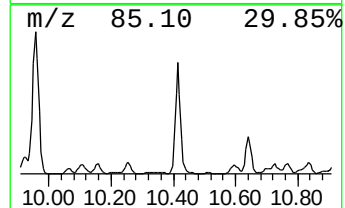
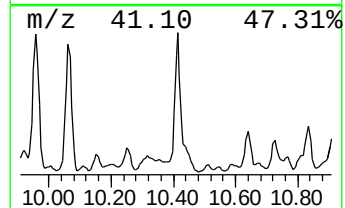
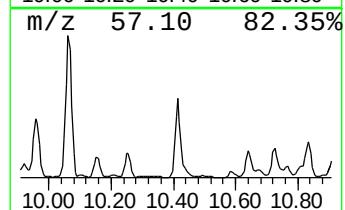
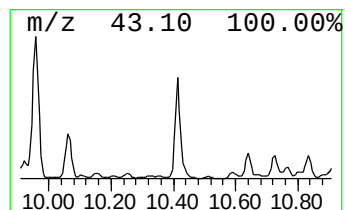
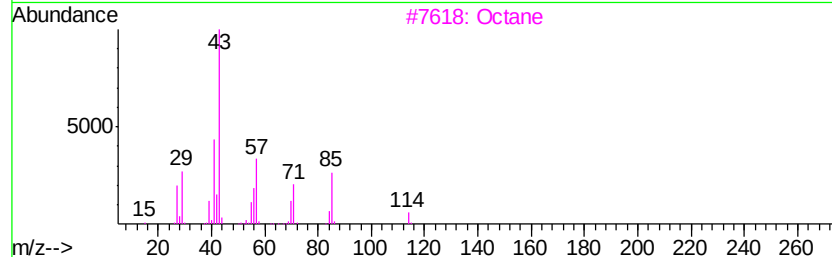
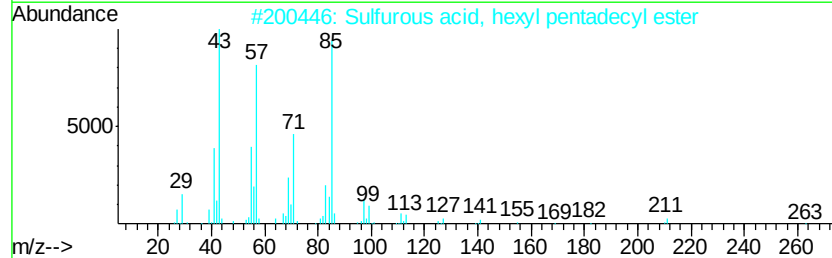
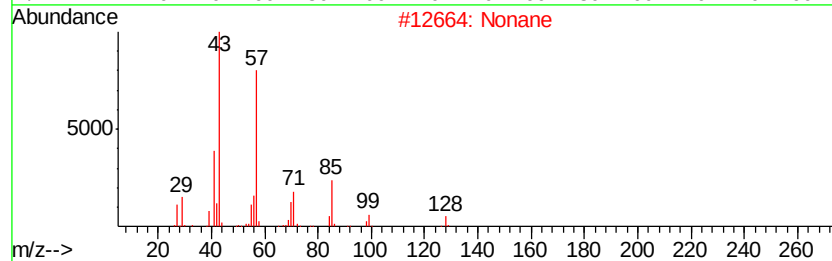
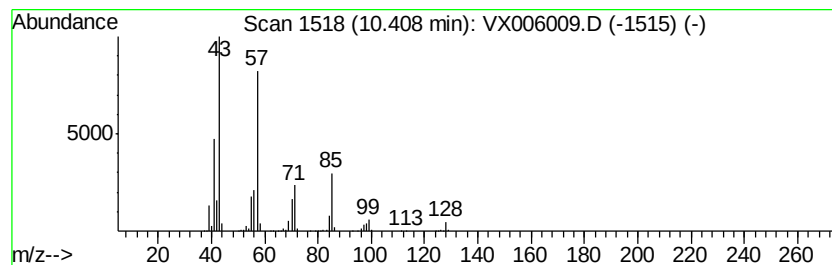
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Peak Number 3 Nonane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.41	54.07 ug/l	549520	Chlorobenzene-d5	10.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane		128	C9H20	000111-84-2	94
2	Sulfurous acid, hexyl pentadecyl...		376	C21H44O3S	1000309-13-7	72
3	Octane		114	C8H18	000111-65-9	72
4	Heptane, 2,4-dimethyl-		128	C9H20	002213-23-2	53
5	Octane, 2,4,6-trimethyl-		156	C11H24	062016-37-9	50



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006009.D
Acq On : 15 Nov 2018 07:39
Operator : JC/MD
Sample : J5915-07
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 46 Sample Multiplier: 1

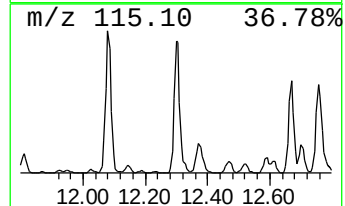
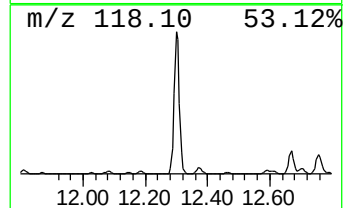
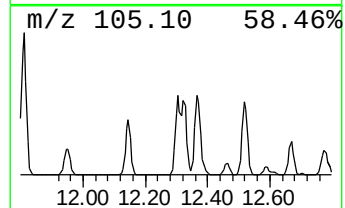
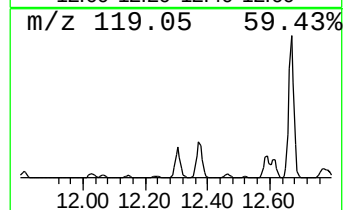
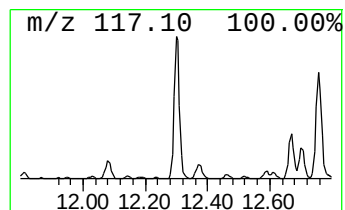
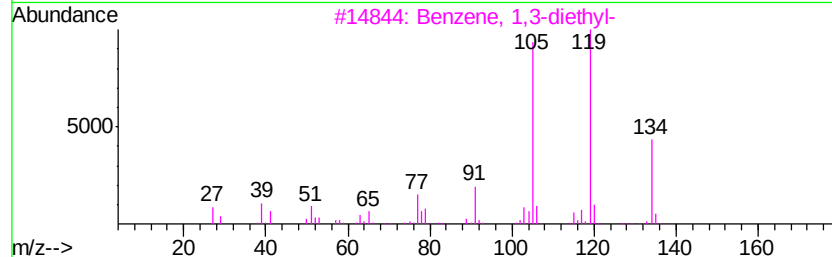
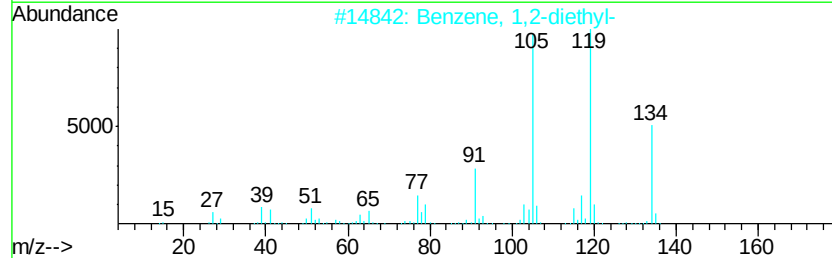
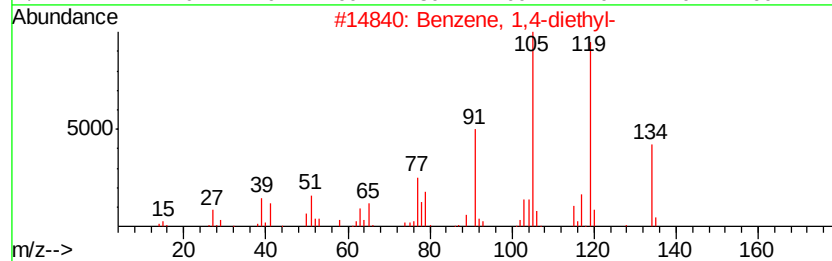
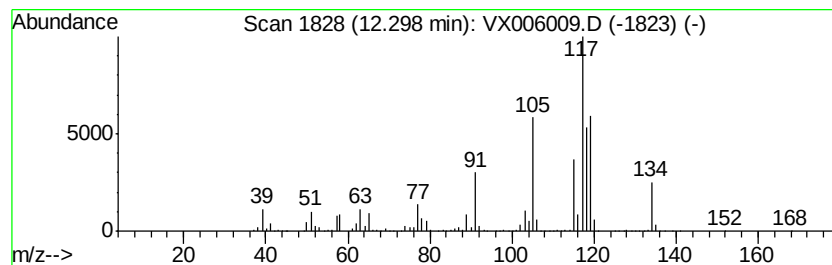
Instrument :
MSVOA_X
ClientSampled :
T-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,4-diethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	78.17 ug/l	1072380	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,4-diethyl-	134 C10H14	000105-05-5 70
2		Benzene, 1,2-diethyl-	134 C10H14	000135-01-3 55
3		Benzene, 1,3-diethyl-	134 C10H14	000141-93-5 55
4		Benzene, cyclopropyl-	118 C9H10	000873-49-4 50
5		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118 C9H10	1000191-13-7 50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006009.D
Acq On : 15 Nov 2018 07:39
Operator : JC/MD
Sample : J5915-07
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 46 Sample Multiplier: 1

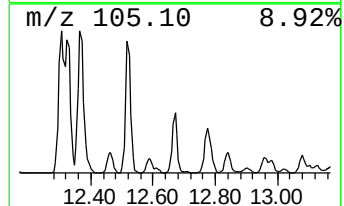
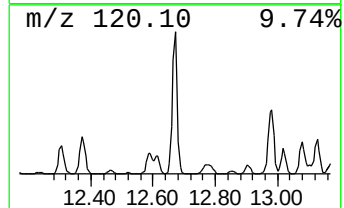
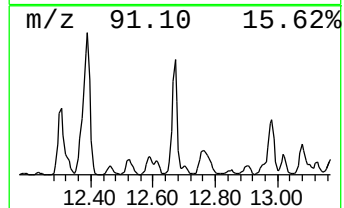
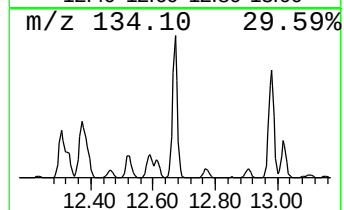
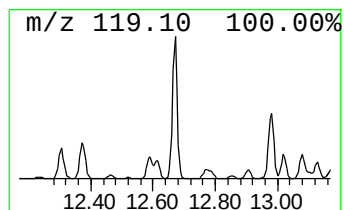
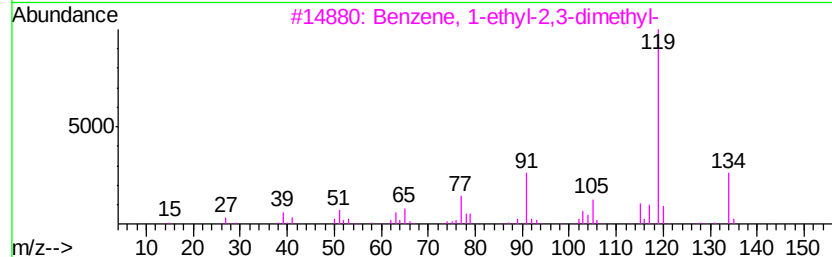
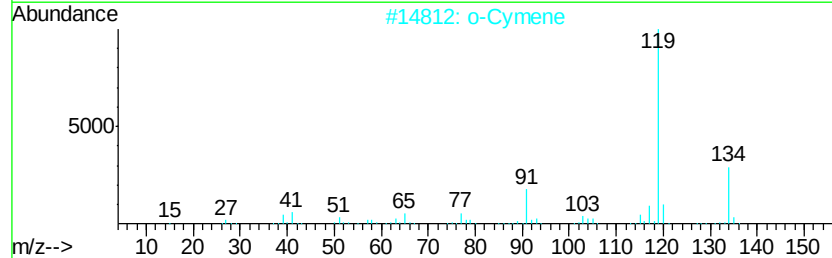
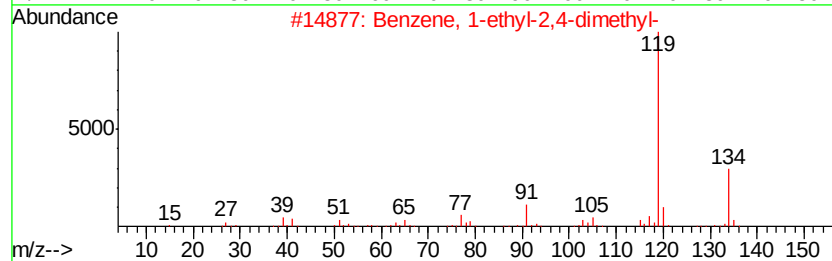
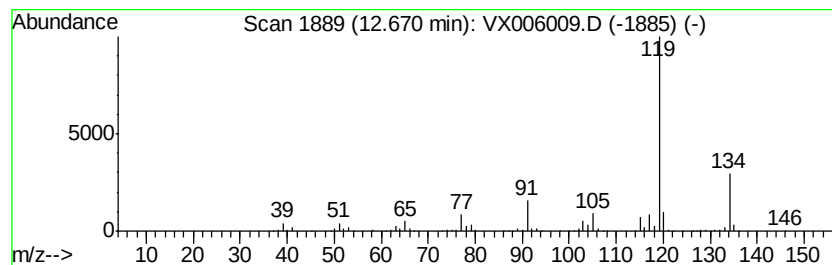
Instrument :
MSVOA_X
ClientSampled :
T-7

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.67	83.65 ug/l	1147610	1,4-Dichlorobenzene-d4	12.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006009.D
Acq On : 15 Nov 2018 07:39
Operator : JC/MD
Sample : J5915-07
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-7

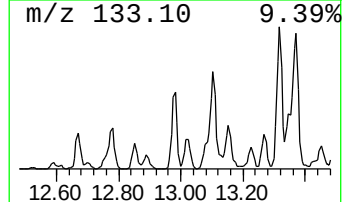
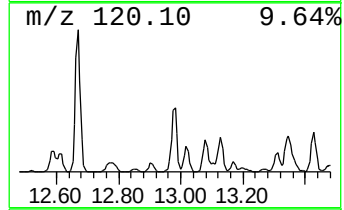
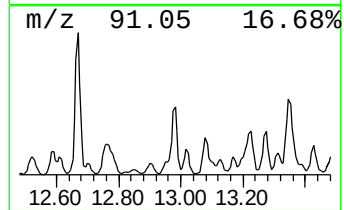
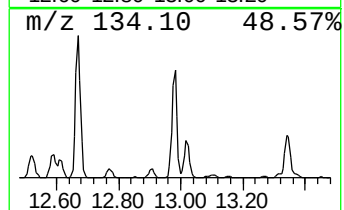
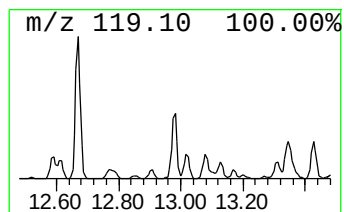
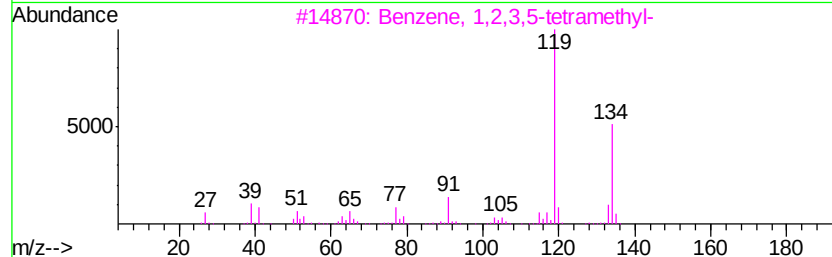
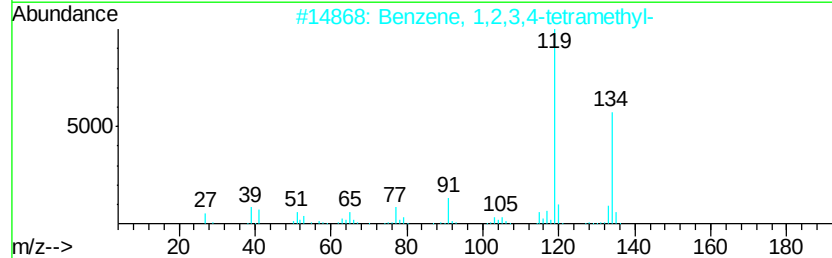
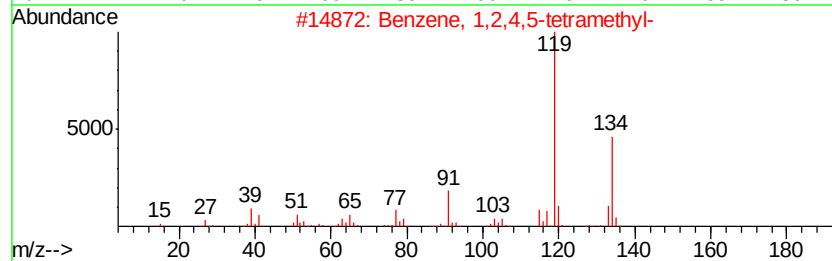
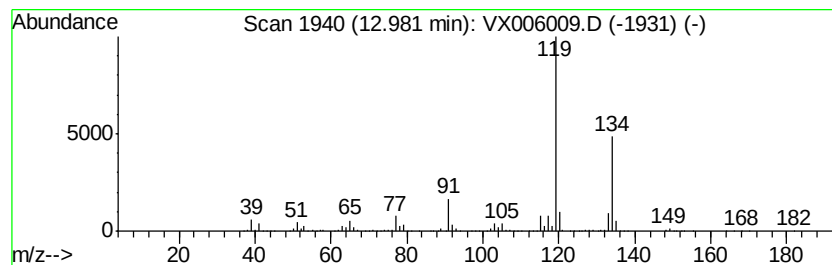
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	60.76 ug/l	833487	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006009.D
Acq On : 15 Nov 2018 07:39
Operator : JC/MD
Sample : J5915-07
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-7

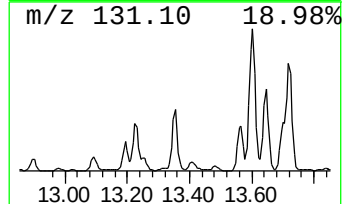
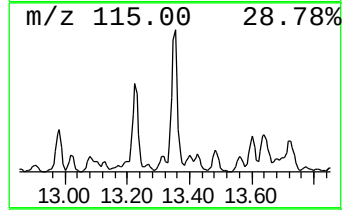
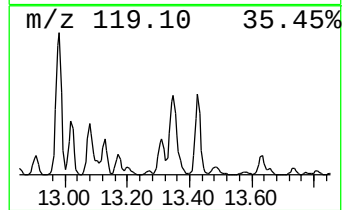
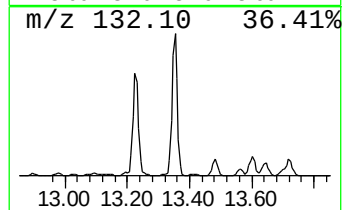
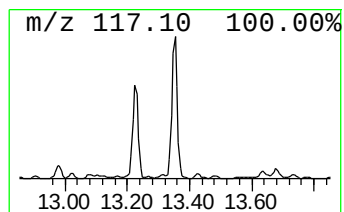
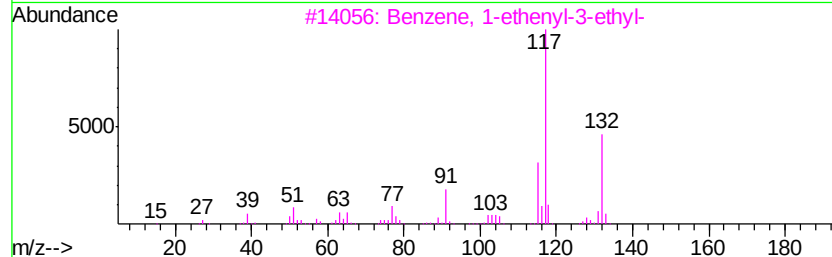
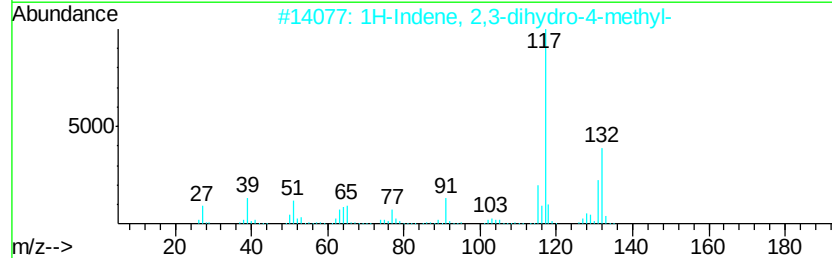
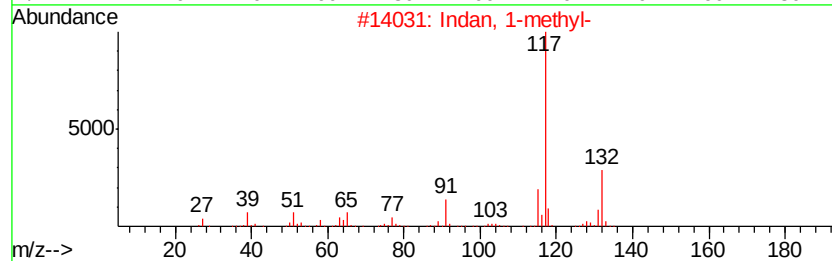
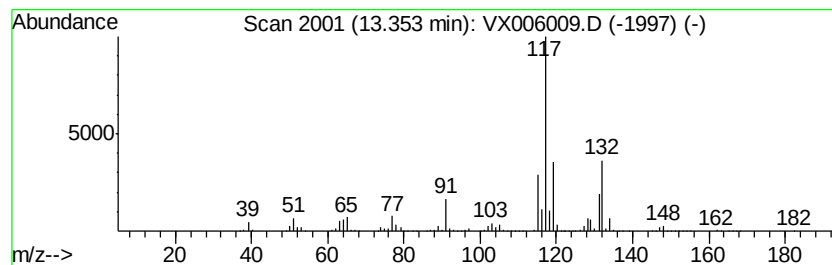
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

Peak Number 9 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	74.46 ug/l	1021490	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Indan, 1-methyl-	132	C10H12	000767-58-8	81
2		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	81
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
5		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	80



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006009.D
Acq On : 15 Nov 2018 07:39
Operator : JC/MD
Sample : J5915-07
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 46 Sample Multiplier: 1

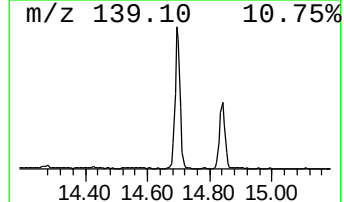
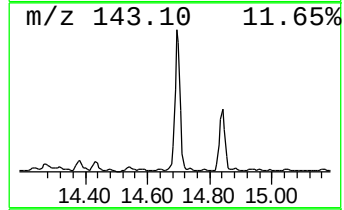
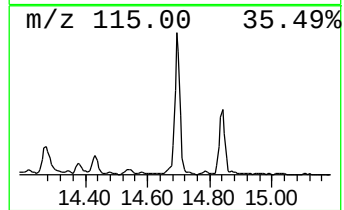
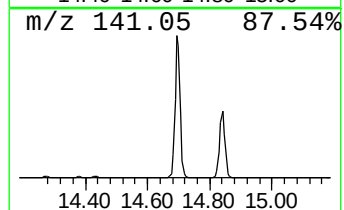
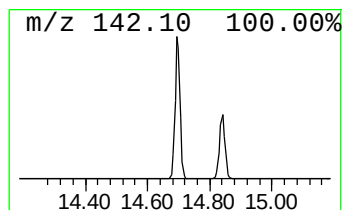
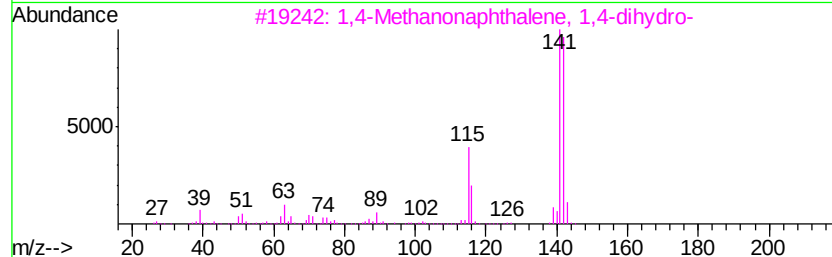
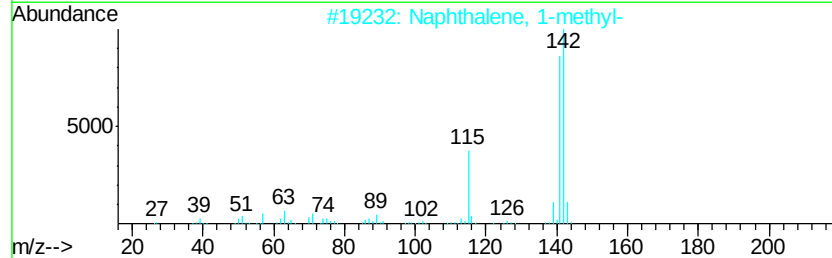
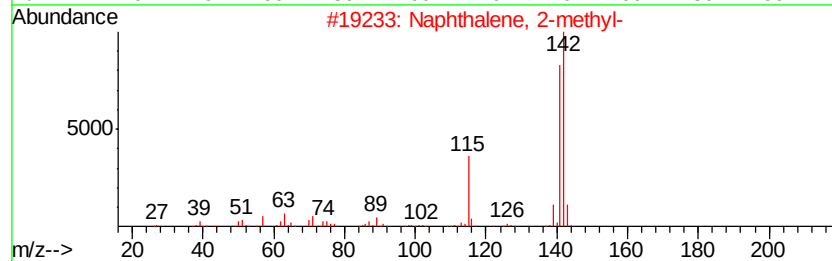
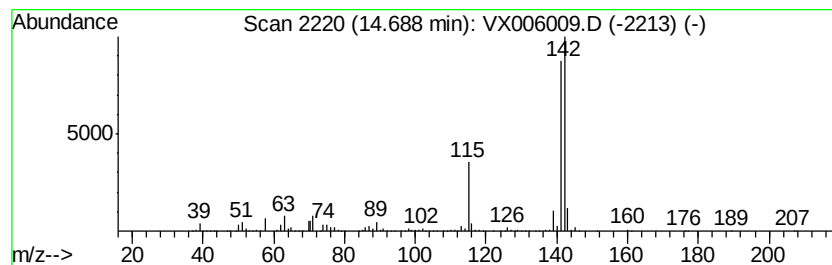
Instrument :
MSVOA_X
ClientSampled :
T-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	45.89 ug/l	629477	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
2		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
3		1,4-Methanonaphthalene, 1,4-dihy...	142 C11H10	004453-90-1 90
4		Benzocycloheptatriene	142 C11H10	000264-09-5 90
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142 C11H10	002443-46-1 45



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006009.D
Acq On : 15 Nov 2018 07:39
Operator : JC/MD
Sample : J5915-07
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-7

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.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
Octane, 4-methyl-	9.96	53.0	ug/l	539049	3	10.12	508165	50.0
Octane, 3-methyl-	10.06	40.9	ug/l	415661	3	10.12	508165	50.0
Nonane	10.41	54.1	ug/l	549520	3	10.12	508165	50.0
Benzene, 1,4-diet...	12.30	78.2	ug/l	1072380	4	12.08	685921	50.0
Benzene, 1-ethyl-...	12.67	83.7	ug/l	1147610	4	12.08	685921	50.0
Benzene, 1,2,4,5-...	12.98	60.8	ug/l	833487	4	12.08	685921	50.0
Indan, 1-methyl-	13.35	74.5	ug/l	1021490	4	12.08	685921	50.0
Naphthalene, 2-me...	14.69	45.9	ug/l	629477	4	12.08	685921	50.0