

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\WX122820\
 Data File : VX020416.D
 Acq On : 28 Dec 2020 16:34
 Operator : JC/MD
 Sample : L5227-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OWL-C4-GW

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121820W.M

Title : Sw846 8260

Signal : TIC: VX020416.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.098	648	658	675	rBV2	18179	57741	2.51%	0.253%
2	5.464	704	718	732	rBV	152101	433917	18.85%	1.901%
3	5.629	732	745	763	rVB	272436	755771	32.82%	3.312%
4	6.031	799	811	831	rBV	172522	450211	19.55%	1.973%
5	6.830	930	942	961	rBV	415659	977956	42.47%	4.285%
6	8.695	1241	1248	1258	rVB	870979	1334625	57.96%	5.848%
7	10.091	1472	1477	1486	rBV	810560	1151548	50.01%	5.046%
8	11.000	1621	1626	1634	rBV	146138	185615	8.06%	0.813%
9	11.122	1640	1646	1652	rBV	679953	872161	37.88%	3.822%
10	11.341	1677	1682	1688	rBV	80427	98878	4.29%	0.433%
11	11.927	1766	1778	1783	rBV	95582	150910	6.55%	0.661%
12	12.061	1795	1800	1807	rBV	886922	1070176	46.48%	4.690%
13	12.213	1820	1825	1831	rVB	40861	52491	2.28%	0.230%
14	12.286	1831	1837	1843	rBV	236263	289287	12.56%	1.268%
15	12.372	1843	1851	1856	rVB	28740	53786	2.34%	0.236%
16	12.445	1856	1863	1868	rBV	92229	117085	5.09%	0.513%
17	12.652	1889	1897	1900	rBV	408799	499839	21.71%	2.190%
18	12.682	1900	1902	1906	rVV	225489	242658	10.54%	1.063%
19	12.737	1906	1911	1919	rVB2	692157	985473	42.80%	4.318%
20	12.878	1926	1934	1939	rBV	49684	74532	3.24%	0.327%
21	12.963	1939	1948	1953	rBV	367952	479468	20.82%	2.101%
22	13.067	1960	1965	1971	rVB2	55944	90939	3.95%	0.398%
23	13.134	1971	1976	1979	rBV	43915	54279	2.36%	0.238%
24	13.176	1979	1983	1985	rVV2	92088	119683	5.20%	0.524%
25	13.207	1985	1988	1998	rVB	302272	390825	16.97%	1.713%
26	13.329	2003	2008	2015	rVV2	1716435	2302497	100.00%	10.090%
27	13.408	2018	2021	2027	rVV2	58708	95853	4.16%	0.420%
28	13.463	2027	2030	2036	rVB	46215	60049	2.61%	0.263%
29	13.542	2039	2043	2046	rBV	60809	76774	3.33%	0.336%
30	13.585	2046	2050	2053	rVV	153780	224238	9.74%	0.983%
31	13.621	2053	2056	2060	rVV2	234899	327101	14.21%	1.433%
32	13.682	2060	2066	2067	rVV2	160739	263398	11.44%	1.154%
33	13.701	2067	2069	2078	rVB2	327809	452143	19.64%	1.981%
34	13.792	2078	2084	2087	rBV	31533	43980	1.91%	0.193%
35	13.835	2087	2091	2095	rBV2	57309	73213	3.18%	0.321%
36	13.890	2095	2100	2105	rVB	349218	448795	19.49%	1.967%
37	13.963	2105	2112	2116	rBV2	463779	613502	26.65%	2.688%

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Sampling : 1

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Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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38	14.011	2116	2120	2124	rVV	132144	167932	7.29%	0.736%
39	14.115	2132	2137	2140	rBV	250687	315625	13.71%	1.383%
40	14.146	2140	2142	2145	rVB	27635	25782	1.12%	0.113%
41	14.255	2154	2160	2169	rBV2	324839	681476	29.60%	2.986%
42	14.359	2173	2177	2181	rVV	131781	157010	6.82%	0.688%
43	14.414	2181	2186	2190	rVV2	271460	361643	15.71%	1.585%
44	14.457	2190	2193	2198	rVB	31201	40799	1.77%	0.179%
45	14.518	2198	2203	2208	rBV2	688743	887381	38.54%	3.889%
46	14.560	2208	2210	2216	rVV4	19069	28109	1.22%	0.123%
47	14.652	2217	2225	2232	rVV3	152758	302997	13.16%	1.328%
48	14.713	2232	2235	2239	rVV	97933	124181	5.39%	0.544%
49	14.767	2239	2244	2250	rVV4	36955	50713	2.20%	0.222%
50	14.828	2250	2254	2256	rVV	43458	67811	2.95%	0.297%
51	14.859	2256	2259	2261	rVV2	73638	103230	4.48%	0.452%
52	14.883	2261	2263	2268	rVV2	53047	68996	3.00%	0.302%
53	14.944	2268	2273	2280	rVV2	58759	120905	5.25%	0.530%
54	15.017	2280	2285	2291	rVB6	11725	26167	1.14%	0.115%
55	15.091	2293	2297	2302	rVB2	27918	38965	1.69%	0.171%
56	15.225	2312	2319	2327	rBV2	153619	304467	13.22%	1.334%
57	15.304	2327	2332	2339	rVB5	39415	72224	3.14%	0.316%
58	15.420	2346	2351	2354	rBV	132001	190128	8.26%	0.833%
59	15.450	2354	2356	2361	rVB	103816	134742	5.85%	0.590%
60	15.523	2362	2368	2373	rVV	438222	688784	29.91%	3.018%
61	15.566	2373	2375	2381	rVB3	50618	79410	3.45%	0.348%
62	15.664	2381	2391	2397	rBV	495850	770941	33.48%	3.378%
63	15.779	2405	2410	2417	rBV3	24846	57110	2.48%	0.250%
64	15.889	2418	2428	2445	rVV	189222	514486	22.34%	2.255%
65	16.042	2447	2453	2465	rVB	115704	207441	9.01%	0.909%
66	16.389	2501	2510	2522	rVB2	44953	124393	5.40%	0.545%
67	16.590	2538	2543	2546	rVV3	10834	23274	1.01%	0.102%
68	16.657	2550	2554	2560	rVV	31251	61123	2.65%	0.268%
69	16.718	2560	2564	2575	rVB3	24012	50725	2.20%	0.222%

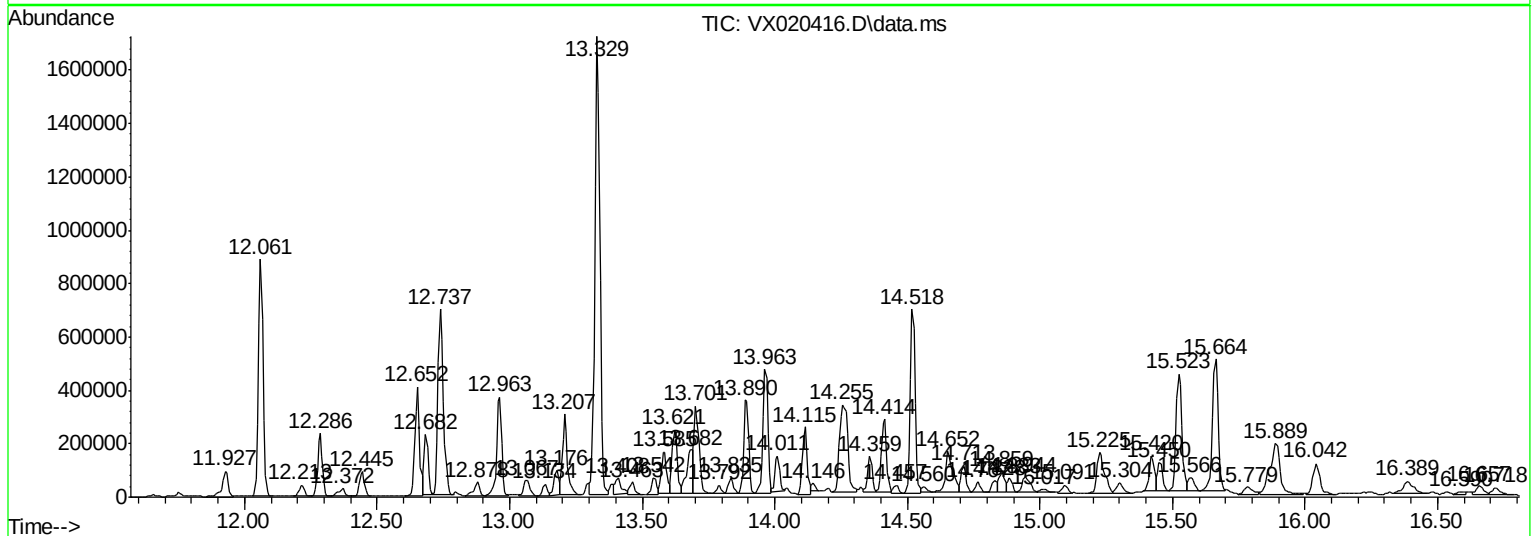
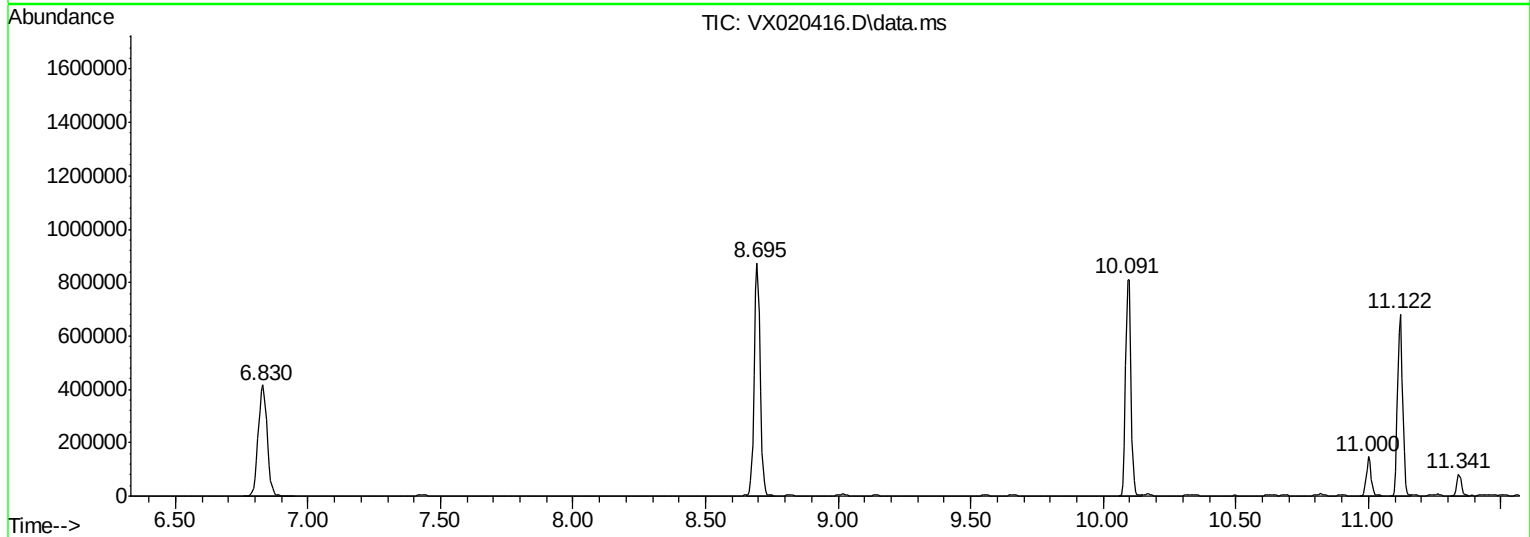
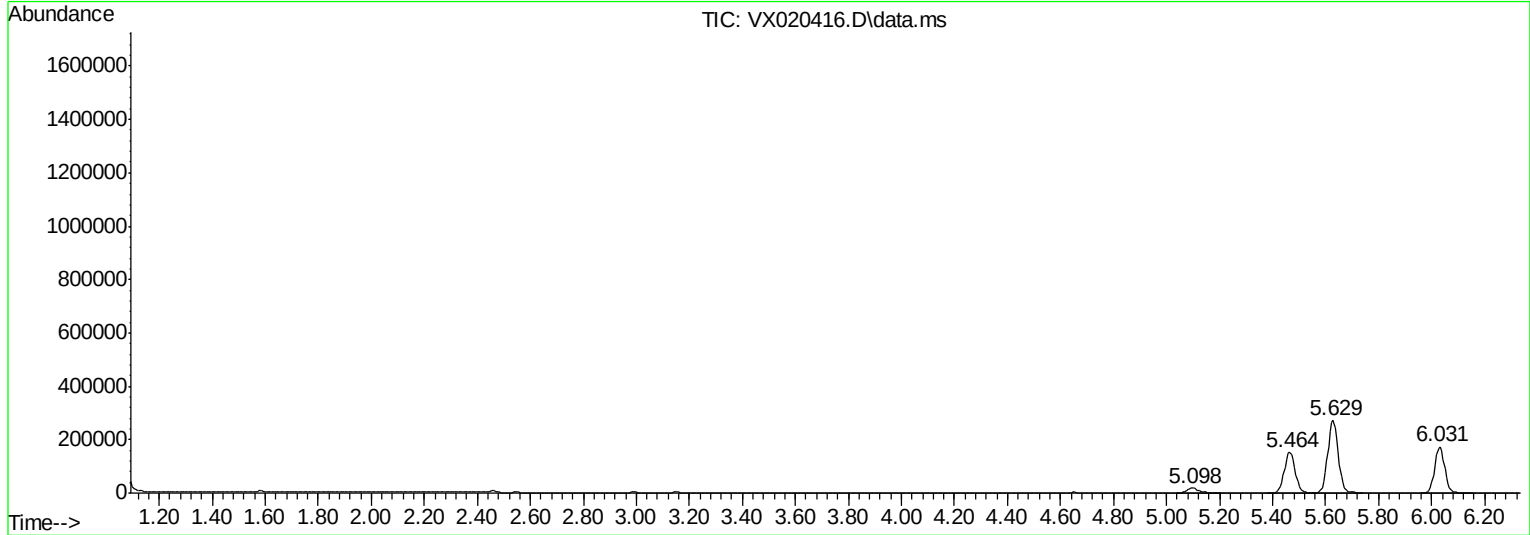
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Instrument :
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ClientSampleId :
OWL-C4-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121820W.M
Quant Title : SW846 8260

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



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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OWL-C4-GW

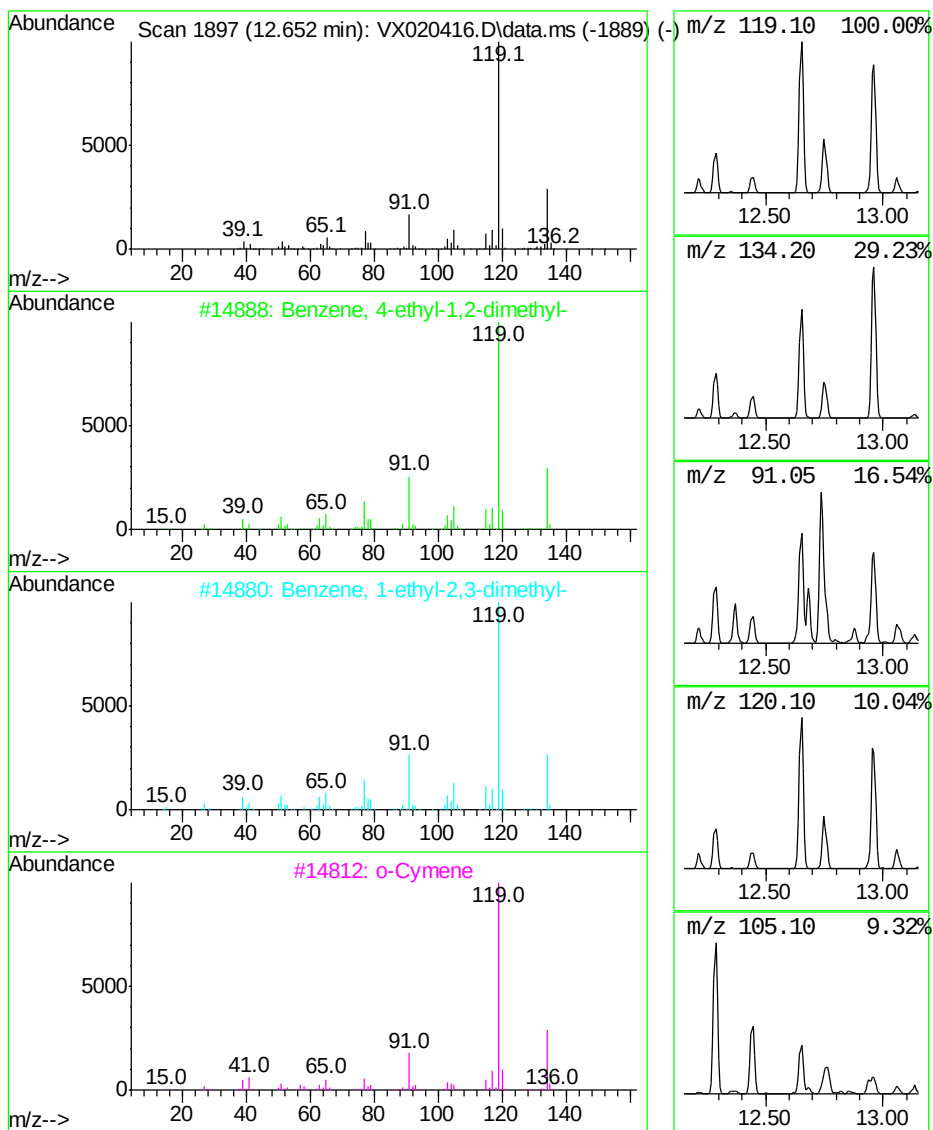
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Quant Title : SW846 8260

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

Peak Number 1 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.652	23.35 ug/l	499839	1,4-Dichlorobenzene-d4	12.061

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



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Sample : L5227-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OWL-C4-GW

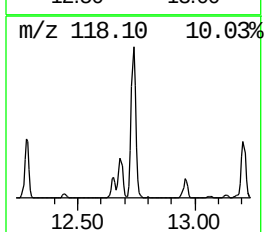
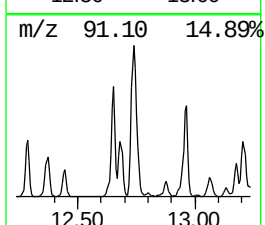
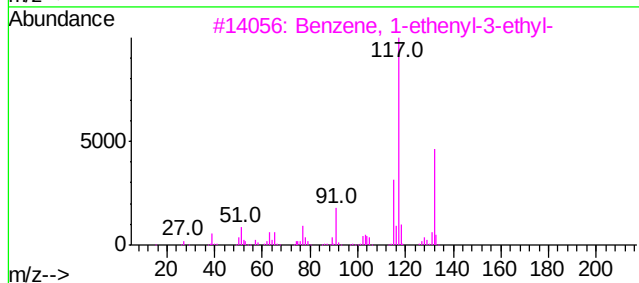
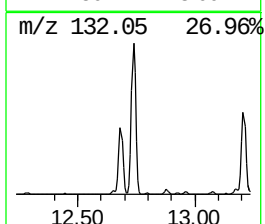
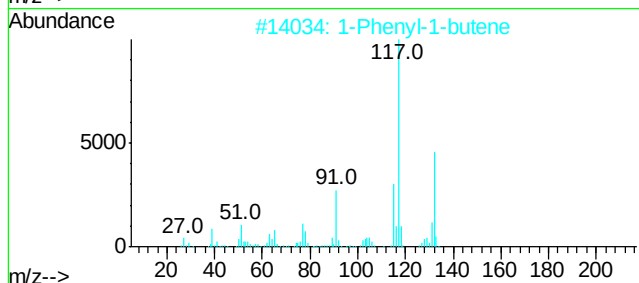
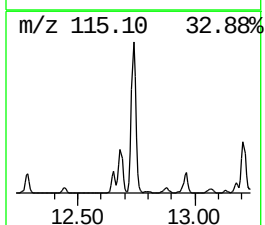
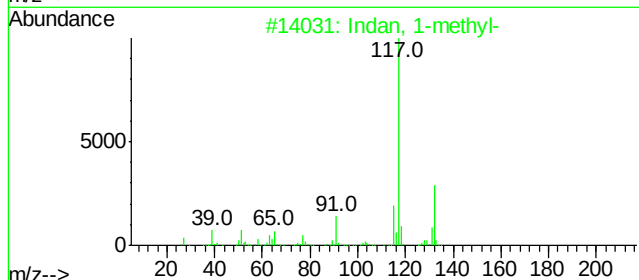
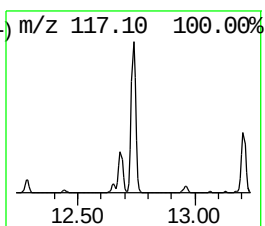
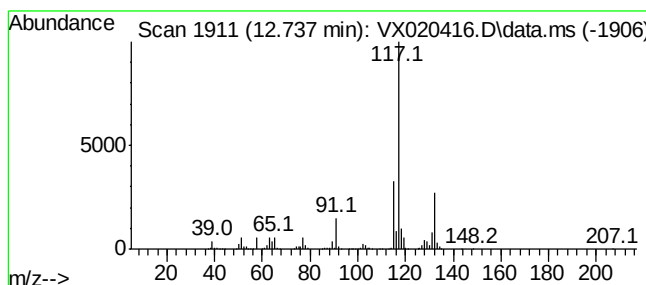
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121820W.M
Quant Title : SW846 8260

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

Peak Number 2 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.737	46.04 ug/l	985473	1,4-Dichlorobenzene-d4	12.061

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	86
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	80



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Sample : L5227-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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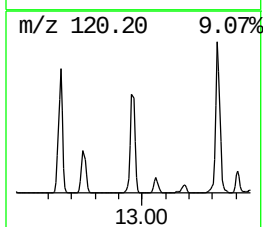
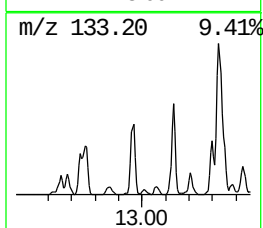
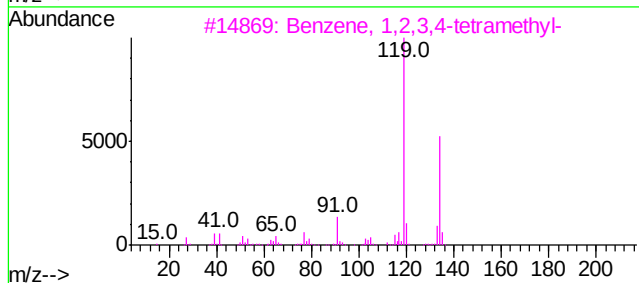
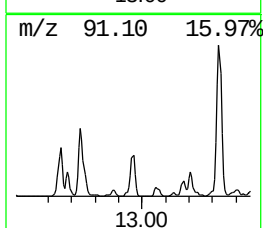
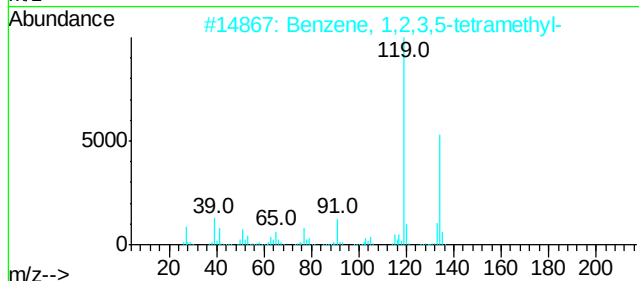
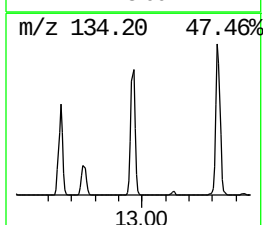
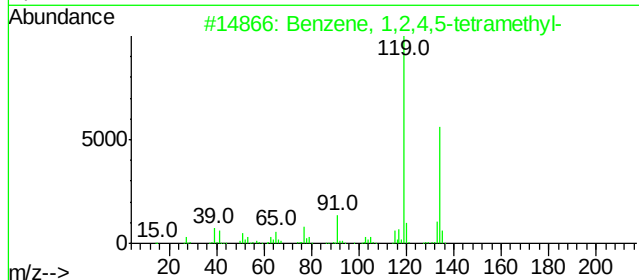
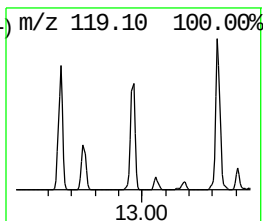
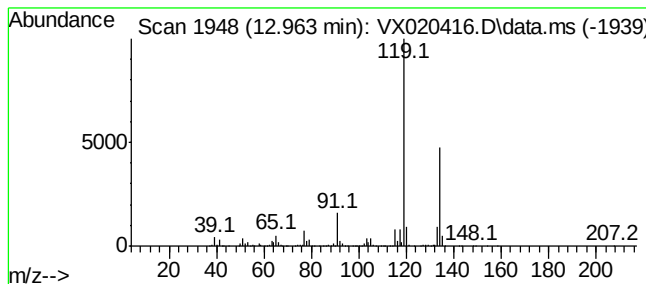
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	22.40 ug/l	479468	1,4-Dichlorobenzene-d4	12.061

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



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Instrument :
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ClientSampleId :
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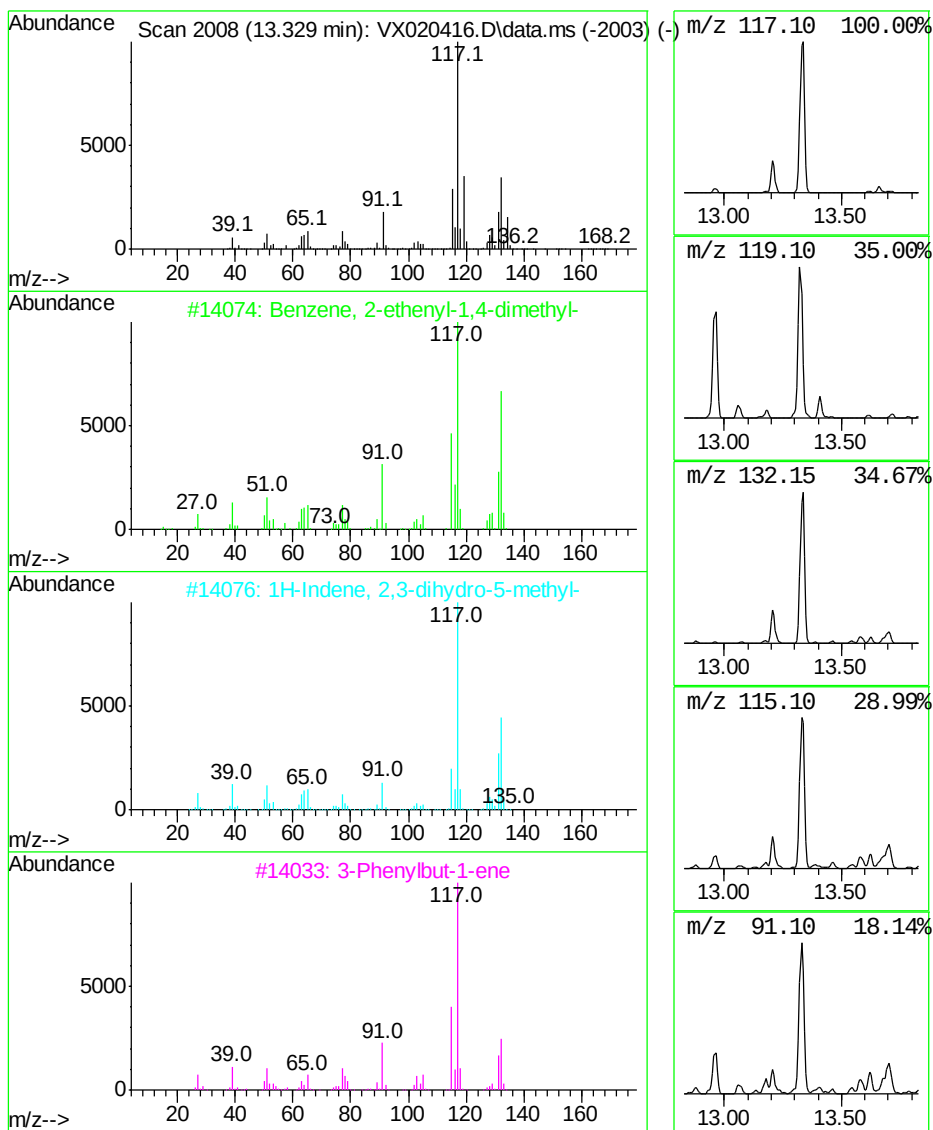
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121820W.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.329	107.58 ug/l	2302500	1,4-Dichlorobenzene-d4	12.061

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83
5	Indan, 1-methyl-	132	C10H12	000767-58-8	81



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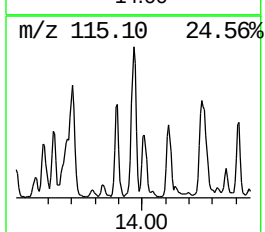
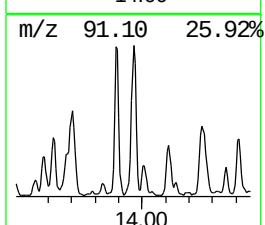
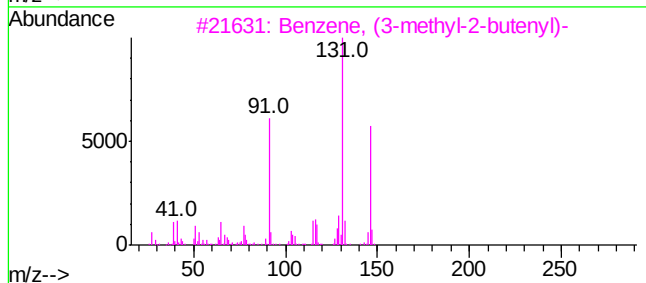
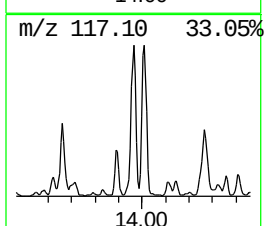
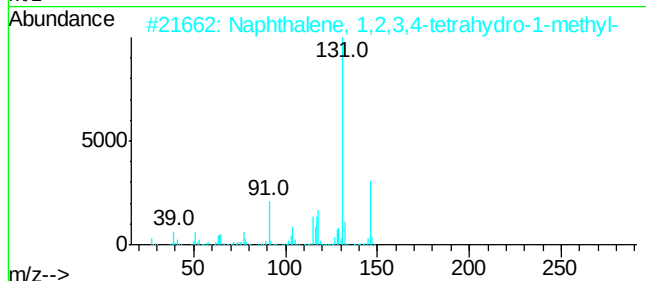
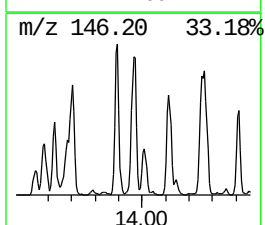
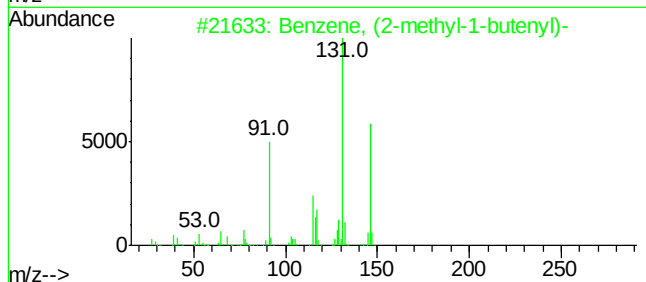
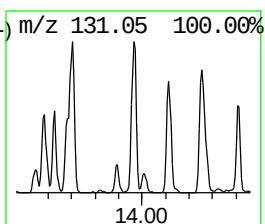
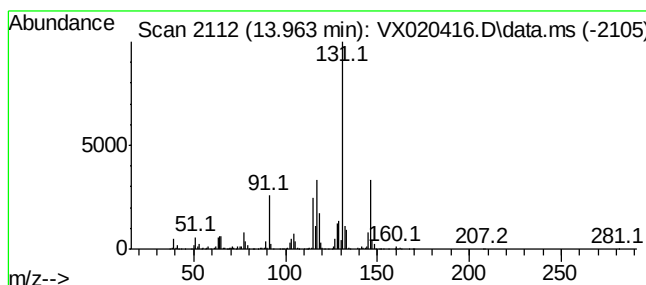
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Peak Number 5 Benzene, (2-methyl-1-butenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.963	28.66 ug/l	613502	1,4-Dichlorobenzene-d4	12.061

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	93
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	93
4	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	87
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122820\
Data File : VX020416.D
Acq On : 28 Dec 2020 16:34
Operator : JC/MD
Sample : L5227-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OWL-C4-GW

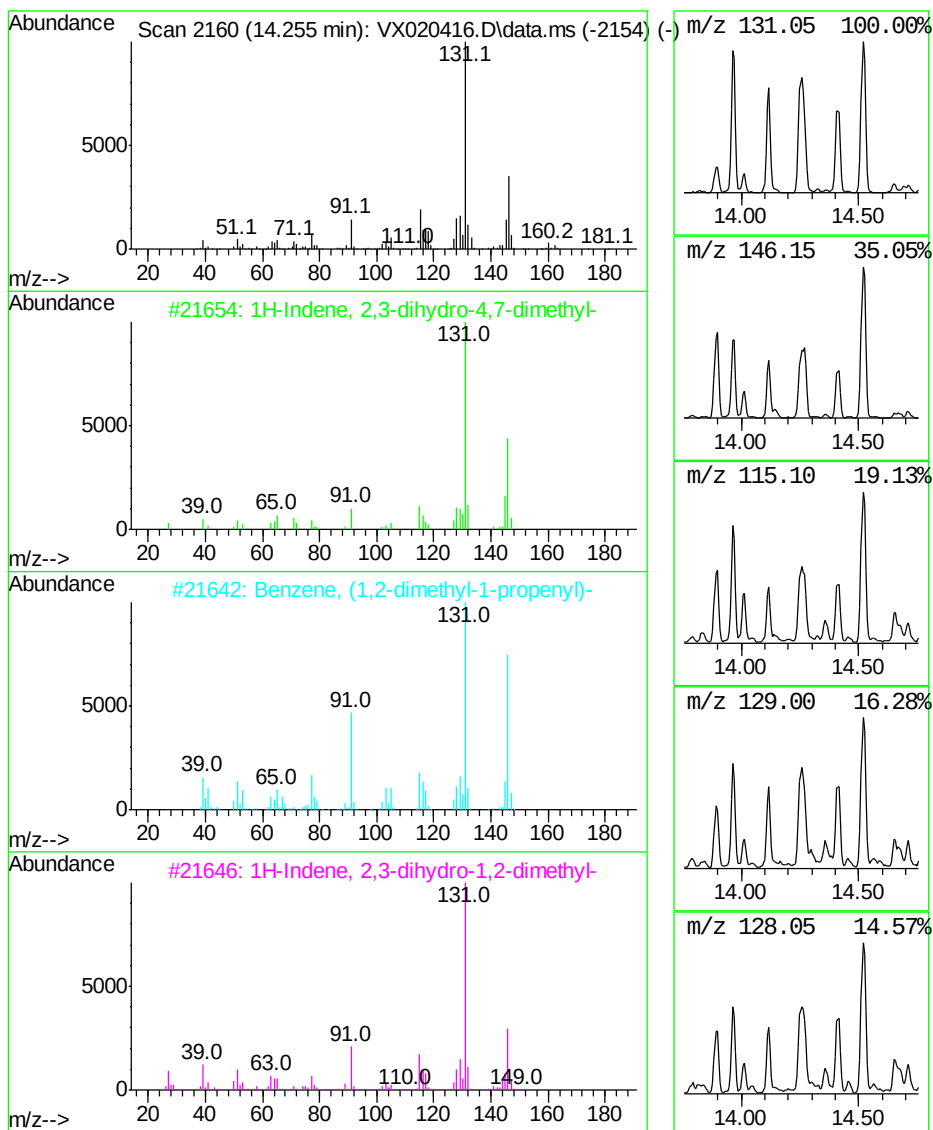
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Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.255	31.84 ug/l	681476	1,4-Dichlorobenzene-d4	12.061

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	96
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	006769-57-3	94
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	94
4	Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	91
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122820\
Data File : VX020416.D
Acq On : 28 Dec 2020 16:34
Operator : JC/MD
Sample : L5227-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OWL-C4-GW

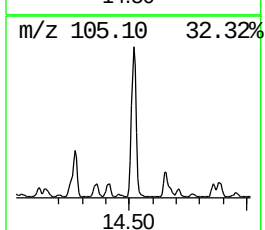
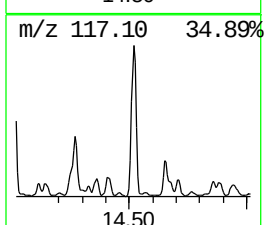
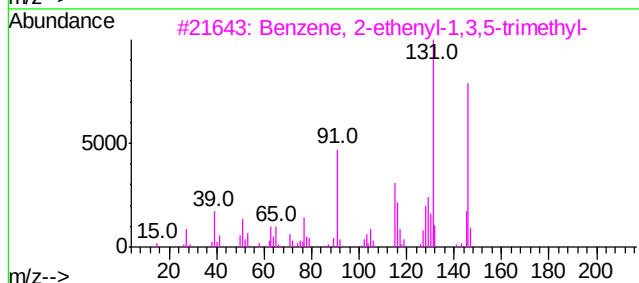
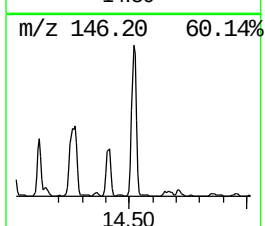
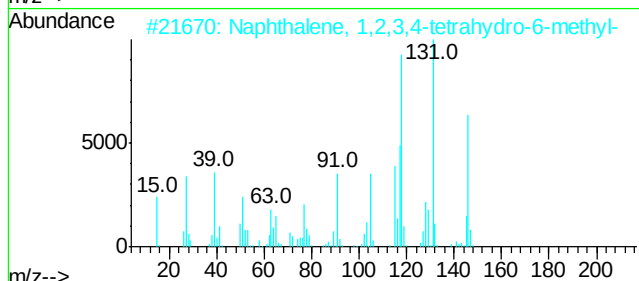
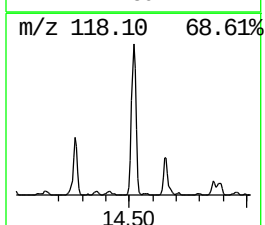
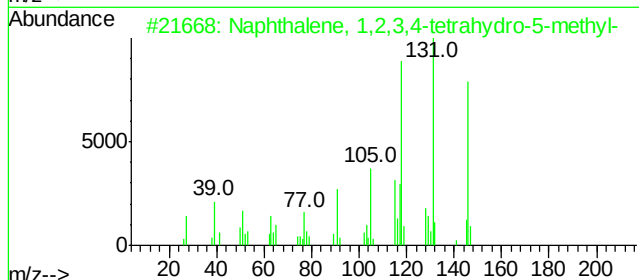
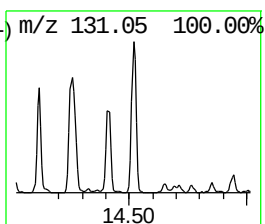
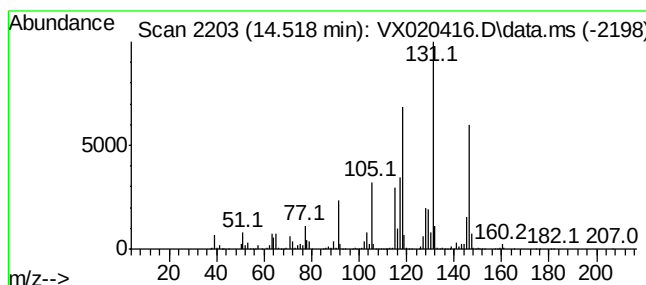
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.518	41.46 ug/l	887381	1,4-Dichlorobenzene-d4	12.061

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
3	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	89
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122820\
Data File : VX020416.D
Acq On : 28 Dec 2020 16:34
Operator : JC/MD
Sample : L5227-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OWL-C4-GW

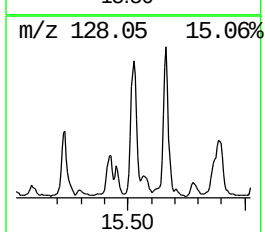
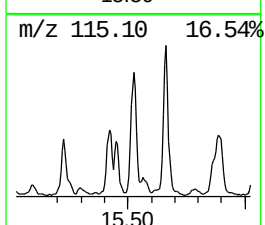
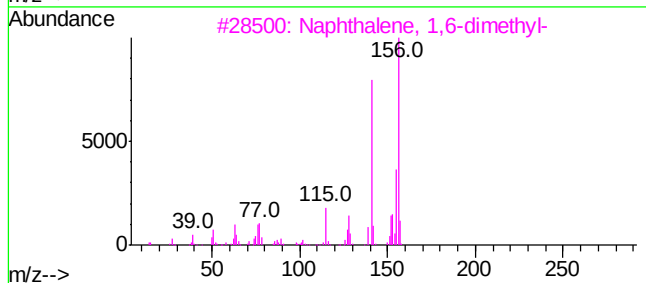
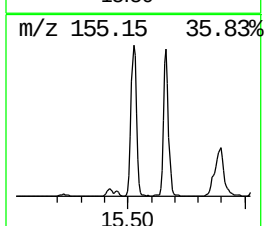
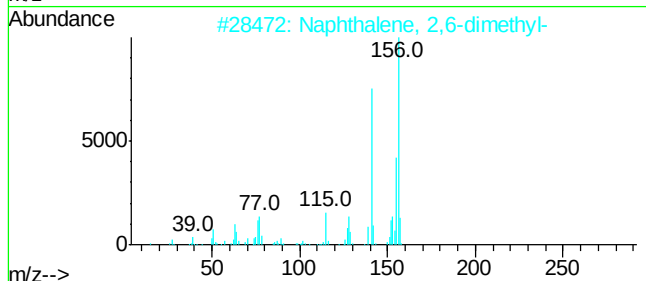
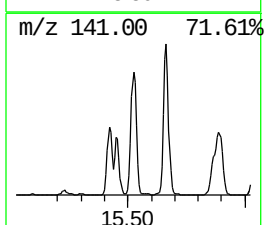
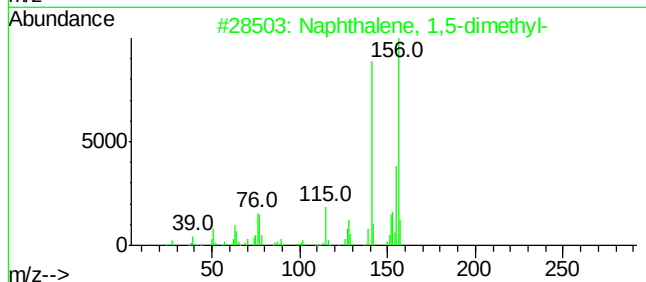
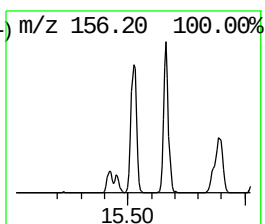
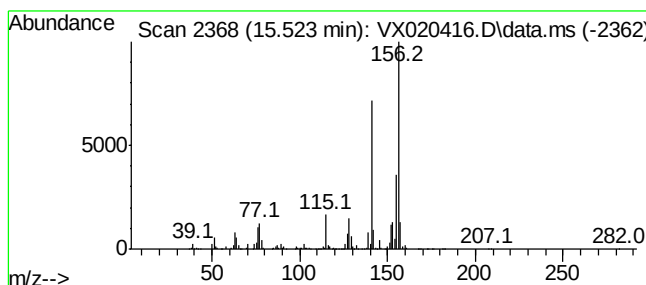
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121820W.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,5-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.523	32.18 ug/l	688784	1,4-Dichlorobenzene-d4	12.061

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	98
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
5	Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122820\
Data File : VX020416.D
Acq On : 28 Dec 2020 16:34
Operator : JC/MD
Sample : L5227-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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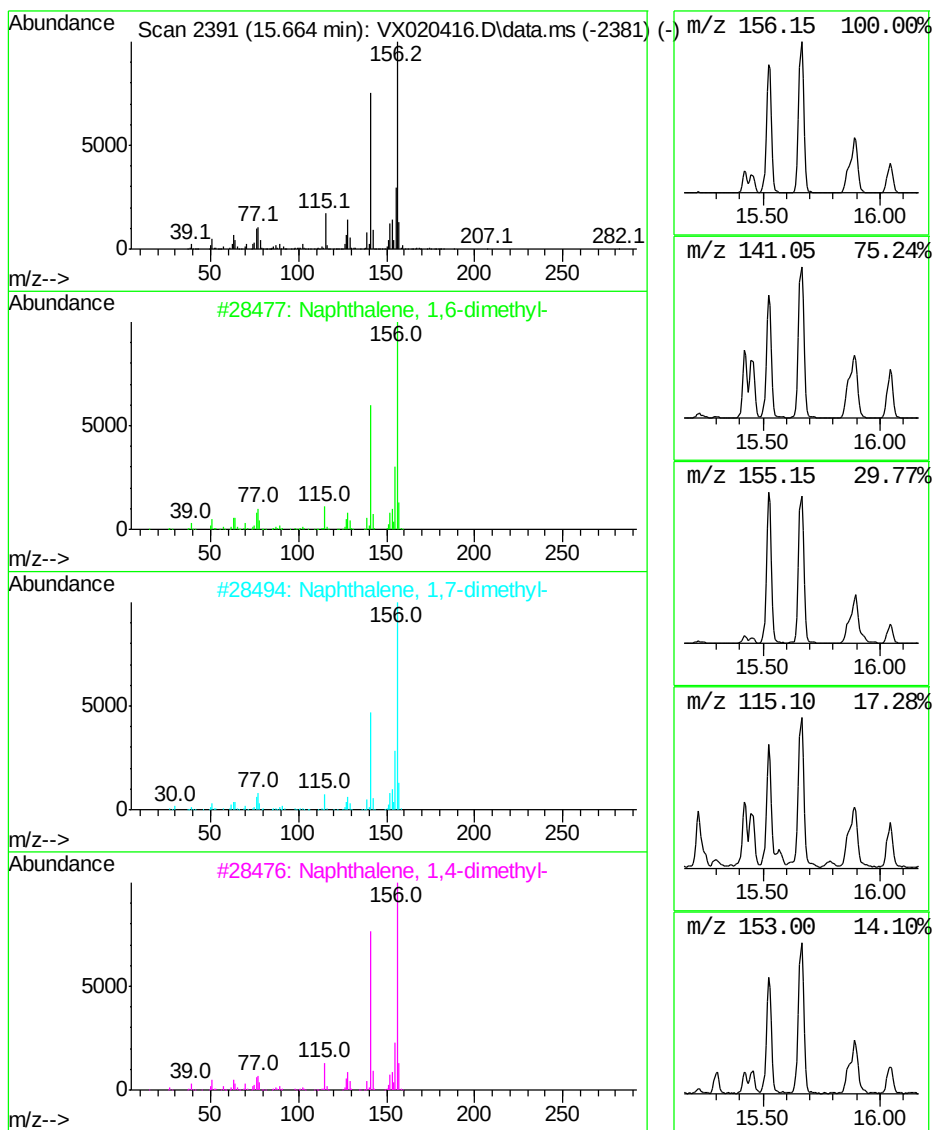
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.664	36.02 ug/l	770941	1,4-Dichlorobenzene-d4	12.061

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122820\
Data File : VX020416.D
Acq On : 28 Dec 2020 16:34
Operator : JC/MD
Sample : L5227-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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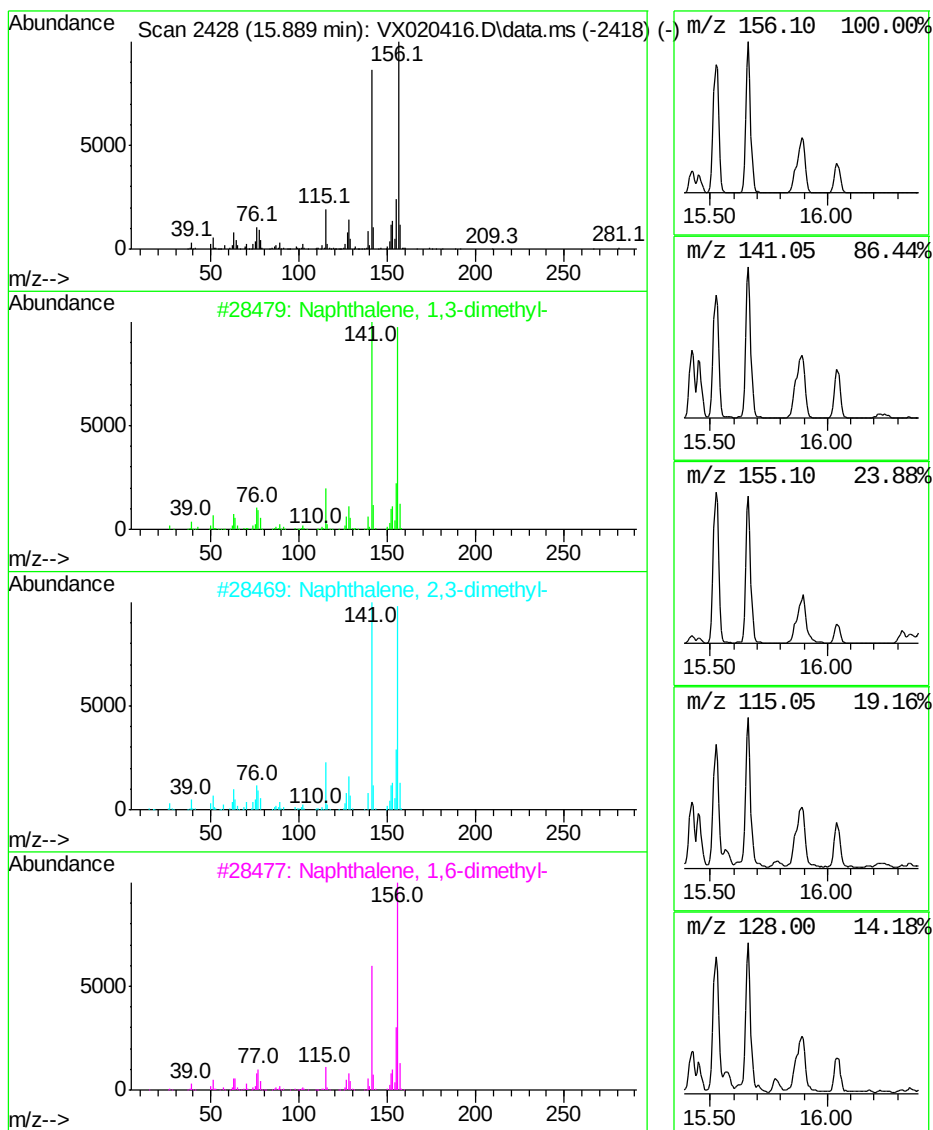
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.889	24.04 ug/l	514486	1,4-Dichlorobenzene-d4	12.061

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	98
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
5	Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122820\
Data File : VX020416.D
Acq On : 28 Dec 2020 16:34
Operator : JC/MD
Sample : L5227-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OWL-C4-GW

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121820W.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
Benzene, 4-ethy...	12.652	23.4	ug/l	499839	4	12.061	1070180	50.0
Indan, 1-methyl-	12.737	46.0	ug/l	985473	4	12.061	1070180	50.0
Benzene, 1,2,4,...	12.963	22.4	ug/l	479468	4	12.061	1070180	50.0
Benzene, 2-ethe...	13.329	107.6	ug/l	2302500	4	12.061	1070180	50.0
Benzene, (2-met...	13.963	28.7	ug/l	613502	4	12.061	1070180	50.0
1H-Indene, 2,3-...	14.255	31.8	ug/l	681476	4	12.061	1070180	50.0
Naphthalene, 1,...	14.518	41.5	ug/l	887381	4	12.061	1070180	50.0
Naphthalene, 1,...	15.523	32.2	ug/l	688784	4	12.061	1070180	50.0
Naphthalene, 1,...	15.664	36.0	ug/l	770941	4	12.061	1070180	50.0
Naphthalene, 1,...	15.889	24.0	ug/l	514486	4	12.061	1070180	50.0