

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012621\
 Data File : VX020765.D
 Acq On : 26 Jan 2021 19:48
 Operator : JC/SP
 Sample : M1217-16 50X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 1710

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\82X012521W.M

Title : SW846 8260

Signal : TIC: VX020765.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.471	60	63	70	rVB	23136	28665	1.48%	0.132%
2	1.587	73	82	87	rBV7	9384	21674	1.12%	0.100%
3	2.331	198	204	211	rBV	9522	21898	1.13%	0.101%
4	2.989	304	312	322	rVB2	8418	21556	1.12%	0.099%
5	5.464	706	718	733	rBV	224818	637931	33.02%	2.935%
6	5.629	733	745	758	rVB	379867	1051178	54.41%	4.837%
7	6.031	800	811	829	rBV2	240489	631519	32.69%	2.906%
8	6.830	928	942	958	rBV	582518	1386302	71.75%	6.379%
9	8.696	1241	1248	1256	rBV	1244797	1932133	100.00%	8.890%
10	10.092	1471	1477	1492	rVB	1267068	1716396	88.83%	7.898%
11	10.342	1511	1518	1524	rBV	41638	67671	3.50%	0.311%
12	10.622	1556	1564	1568	rBV5	10959	21488	1.11%	0.099%
13	10.683	1568	1574	1581	rVB	25323	40766	2.11%	0.188%
14	10.896	1604	1609	1613	rBV3	18067	28851	1.49%	0.133%
15	10.982	1618	1623	1629	rBV4	11364	20474	1.06%	0.094%
16	11.122	1640	1646	1656	rBV	1034396	1344365	69.58%	6.186%
17	11.415	1690	1694	1701	rBV	49030	97695	5.06%	0.450%
18	11.488	1701	1706	1708	rVV3	32361	54814	2.84%	0.252%
19	11.512	1708	1710	1715	rVB3	27878	34698	1.80%	0.160%
20	11.652	1728	1733	1740	rBV3	35519	60017	3.11%	0.276%
21	11.793	1751	1756	1761	rBV	123973	163246	8.45%	0.751%
22	11.902	1766	1774	1775	rBV5	12968	24981	1.29%	0.115%
23	11.982	1782	1787	1789	rBV4	11894	21742	1.13%	0.100%
24	12.061	1794	1800	1806	rVB	1387570	1695139	87.73%	7.800%
25	12.140	1806	1813	1827	rBV2	1122072	1492058	77.22%	6.865%
26	12.280	1833	1836	1838	rBV	24124	30714	1.59%	0.141%
27	12.305	1838	1840	1844	rVB	43856	46329	2.40%	0.213%
28	12.353	1844	1848	1855	rVB4	84978	135073	6.99%	0.622%
29	12.457	1862	1865	1869	rVB3	28099	29460	1.52%	0.136%
30	12.500	1869	1872	1878	rVB3	33552	52830	2.73%	0.243%
31	12.567	1878	1883	1885	rBV2	47998	69142	3.58%	0.318%
32	12.591	1885	1887	1892	rVV	34464	41391	2.14%	0.190%
33	12.646	1894	1896	1900	rVB	34170	39498	2.04%	0.182%
34	12.701	1900	1905	1908	rBV4	23294	36571	1.89%	0.168%
35	12.756	1908	1914	1918	rBV5	43135	94004	4.87%	0.433%
36	12.853	1925	1930	1933	rBV2	104041	148791	7.70%	0.685%

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37	12.890	1933	1936	1939	rVB3	53079	62867	3.25%	0.289%
38	12.963	1939	1948	1951	rBV4	63530	112144	5.80%	0.516%
39	13.006	1951	1955	1963	rBV5	79413	152705	7.90%	0.703%
40	13.079	1964	1967	1969	rBV4	20288	24342	1.26%	0.112%
41	13.109	1969	1972	1973	rVV	41455	45124	2.34%	0.208%
42	13.134	1973	1976	1981	rVB	126031	152515	7.89%	0.702%
43	13.292	1997	2002	2005	rVV4	29819	64275	3.33%	0.296%
44	13.329	2005	2008	2013	rVB2	157594	218116	11.29%	1.004%
45	13.457	2023	2029	2036	rBV2	801810	1189153	61.55%	5.472%
46	13.548	2040	2044	2048	rBV3	176901	264779	13.70%	1.218%
47	13.628	2053	2057	2060	rVB4	48833	71281	3.69%	0.328%
48	13.707	2067	2070	2072	rBV2	34353	45381	2.35%	0.209%
49	13.743	2073	2076	2079	rVV	193895	213851	11.07%	0.984%
50	13.841	2089	2092	2095	rVB2	67422	79195	4.10%	0.364%
51	13.890	2095	2100	2104	rBV2	192737	305434	15.81%	1.405%
52	13.926	2104	2106	2109	rVV	50377	57564	2.98%	0.265%
53	14.109	2132	2136	2139	rBV4	80568	115086	5.96%	0.530%
54	14.231	2148	2156	2158	rBV3	249410	410738	21.26%	1.890%
55	14.268	2159	2162	2166	rVB4	256499	366649	18.98%	1.687%
56	14.445	2187	2191	2194	rBV	269016	306551	15.87%	1.411%
57	14.524	2200	2204	2212	rBV3	217486	347213	17.97%	1.598%
58	14.591	2212	2215	2217	rBV3	56040	70666	3.66%	0.325%
59	14.658	2222	2226	2232	rVV2	266998	435329	22.53%	2.003%
60	14.713	2232	2235	2238	rBV	194850	224985	11.64%	1.035%
61	14.792	2246	2248	2251	rBV2	97960	110289	5.71%	0.507%
62	14.829	2252	2254	2257	rVV3	95428	113890	5.89%	0.524%
63	14.890	2261	2264	2268	rVB4	136550	171498	8.88%	0.789%
64	15.225	2315	2319	2322	rBV3	164238	272121	14.08%	1.252%
65	15.426	2349	2352	2355	rBV2	124220	153618	7.95%	0.707%
66	15.469	2355	2359	2362	rVV3	111520	148717	7.70%	0.684%
67	15.609	2377	2382	2388	rVB	1257587	1794199	92.86%	8.256%
68	16.139	2467	2469	2475	rVB4	95192	151952	7.86%	0.699%
69	16.773	2570	2573	2575	rBV2	113207	140031	7.25%	0.644%

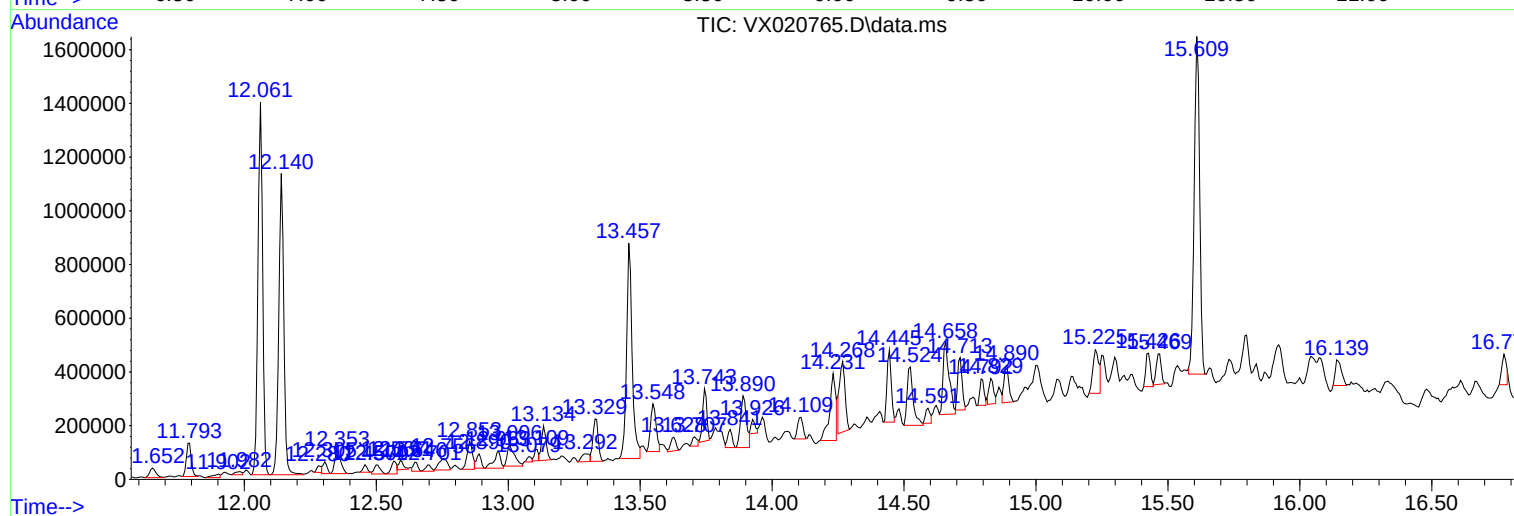
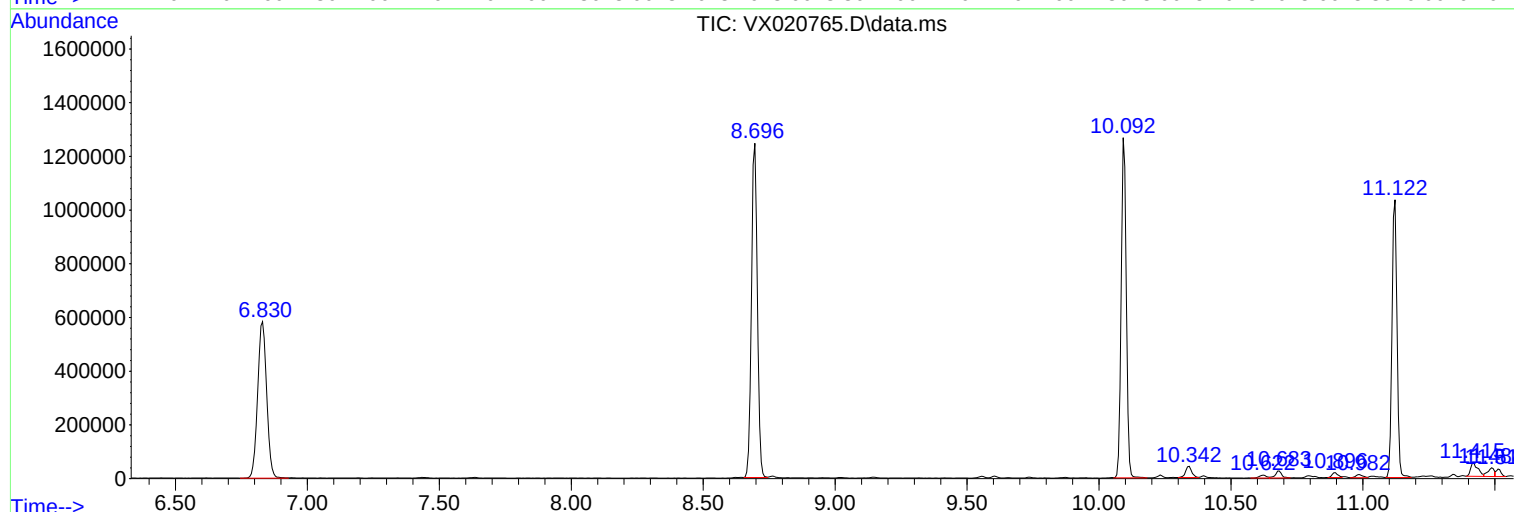
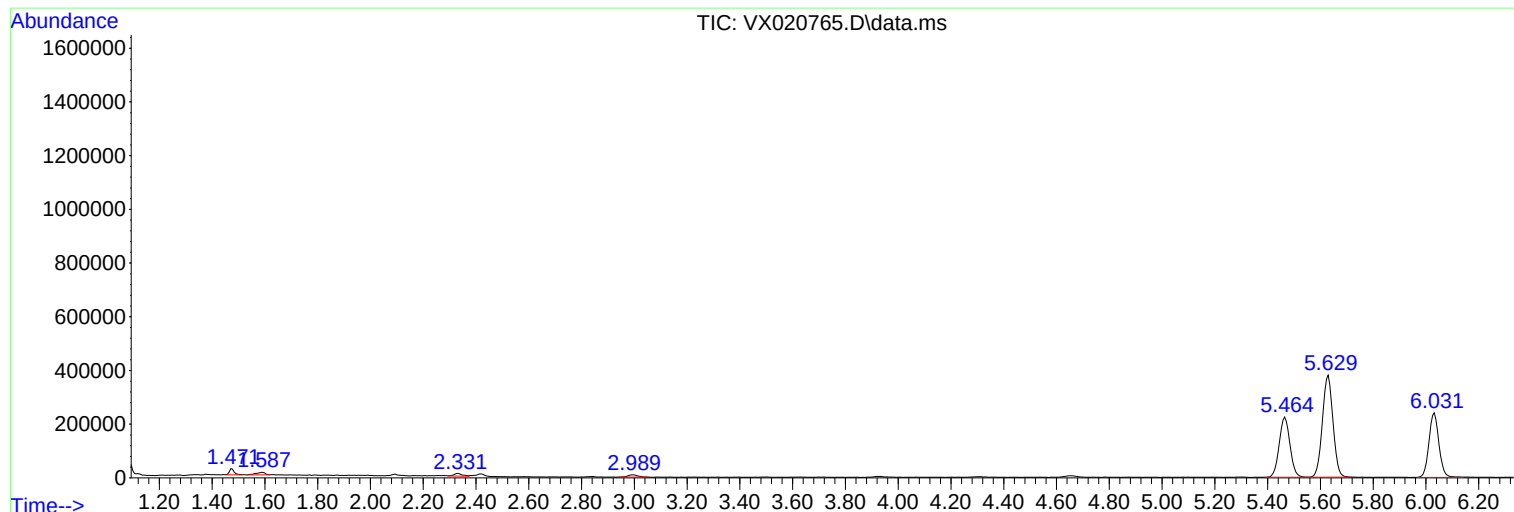
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Instrument :
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ClientSampleId :
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X012521W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_X
ClientSampleId :
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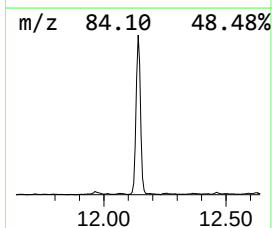
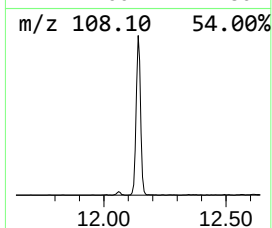
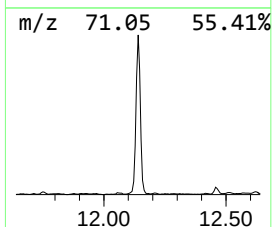
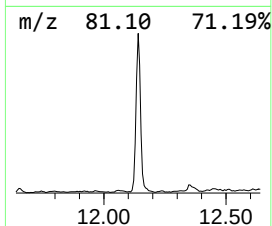
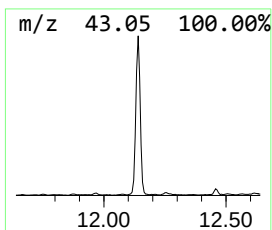
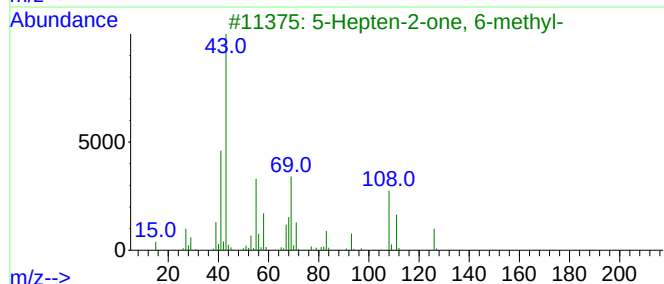
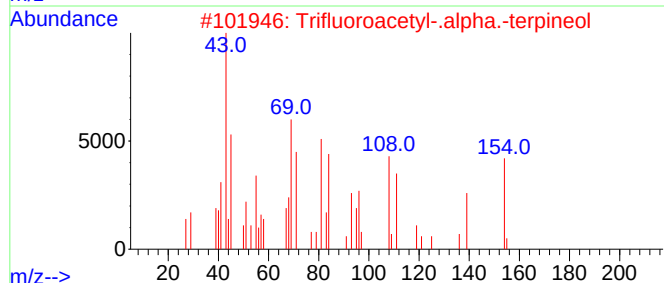
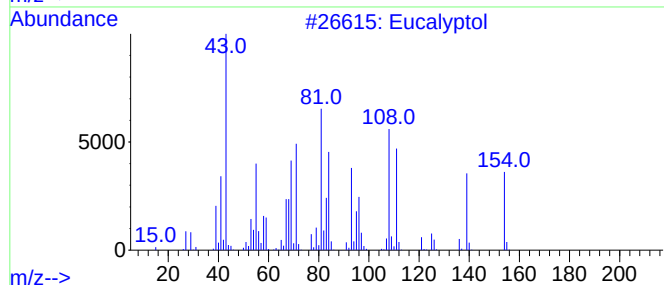
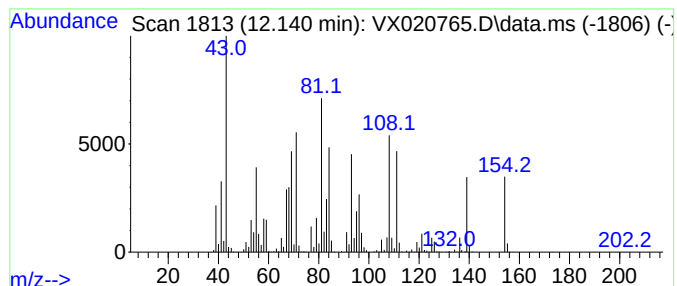
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X012521W.M
Quant Title : SW846 8260

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

Peak Number 1 Eucalyptol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.140	44.01 ug/l	1492060	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eucalyptol	154	C10H18O	000470-82-6	99
2			Trifluoroacetyl-.alpha.-terpineol	250	C12H17F3O2	1000058-17-6	50
3			5-Hepten-2-one, 6-methyl-	126	C8H14O	000110-93-0	35
4			Terpinen-4-ol	154	C10H18O	000562-74-3	25
5			2-Isopropyl-5-methylhex-2-enal	154	C10H18O	035158-25-9	25



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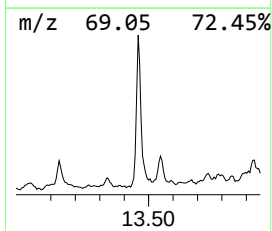
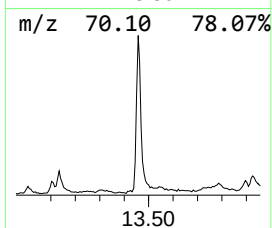
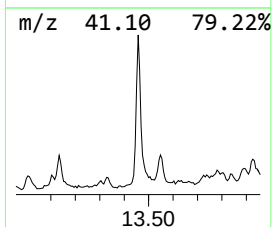
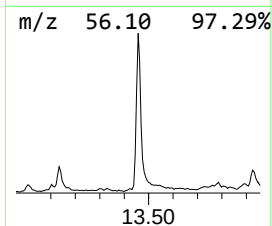
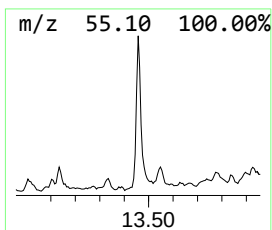
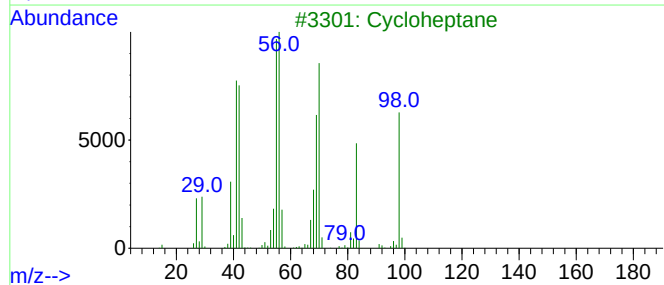
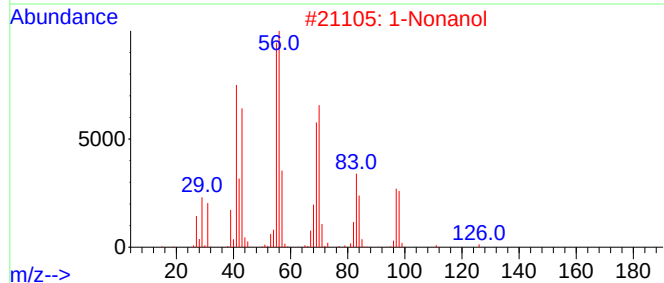
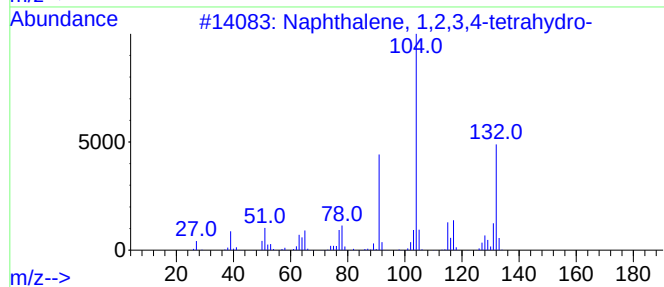
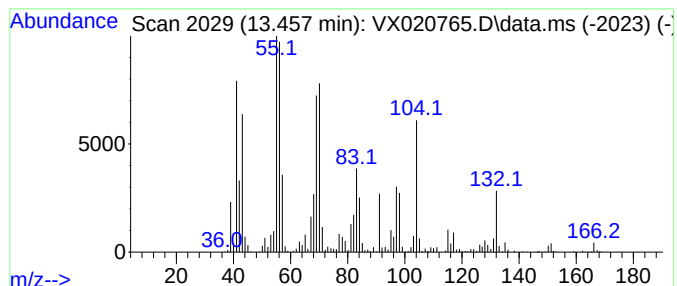
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.457	35.08 ug/l	1189150	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
2			1-Nonanol	144	C9H20O	000143-08-8	68
3			Cycloheptane	98	C7H14	000291-64-5	58
4			Chloroacetic acid, nonyl ester	220	C11H21ClO2	005451-96-7	58
5			1-Decene	140	C10H20	000872-05-9	53



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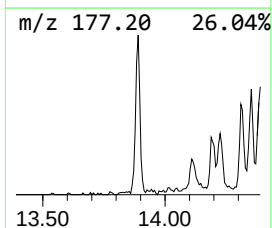
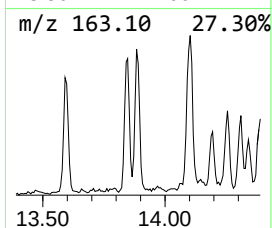
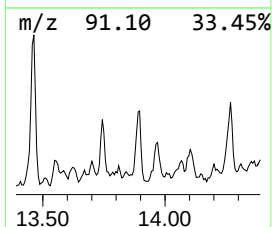
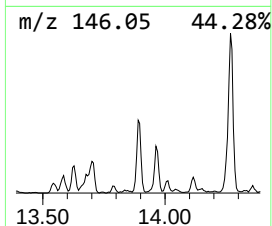
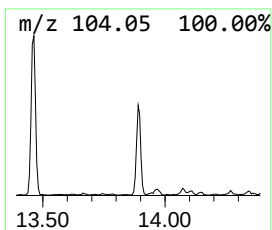
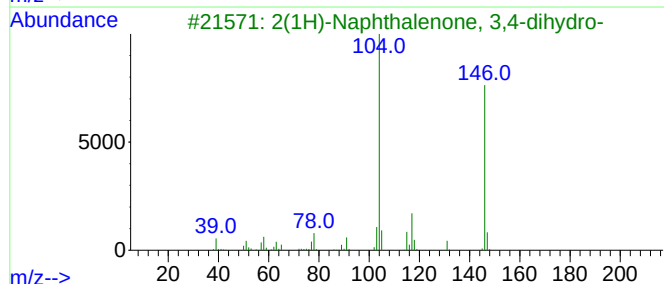
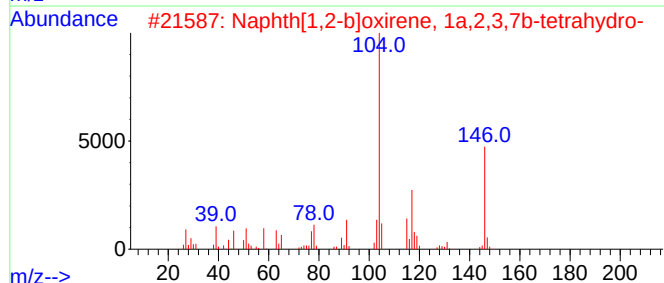
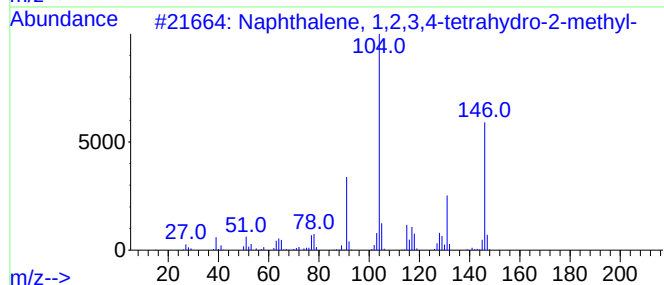
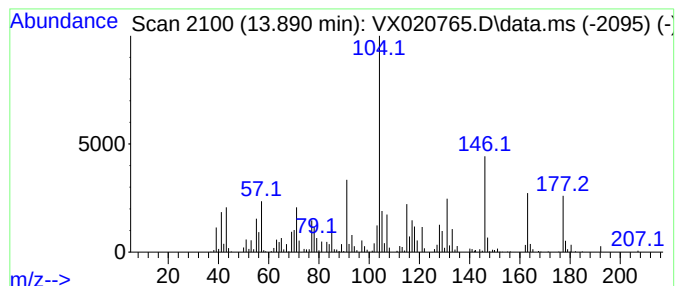
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Peak Number 3 Naphthalene, 1,2,3,4-tetra... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.890	9.01 ug/l	305434	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	86
2			Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	46
3			2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	46
4			Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	38
5			Phthalic acid, methyl 2-phenylet...	284	C17H16O4	1000315-57-6	35



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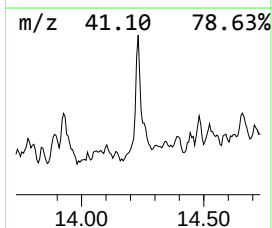
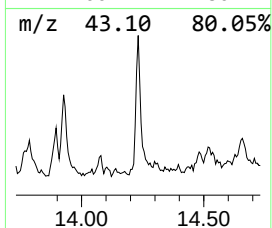
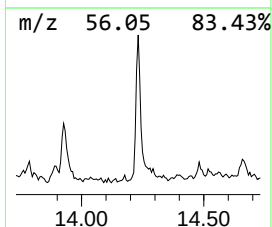
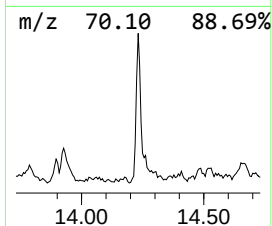
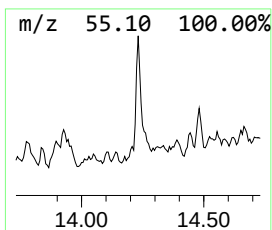
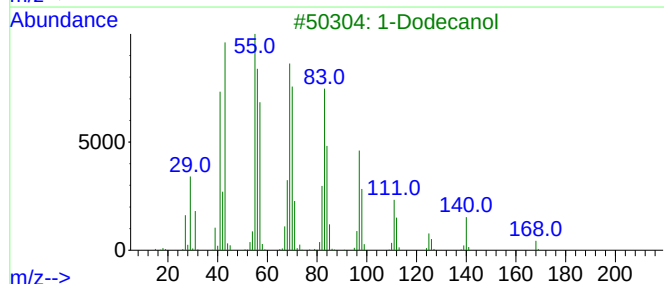
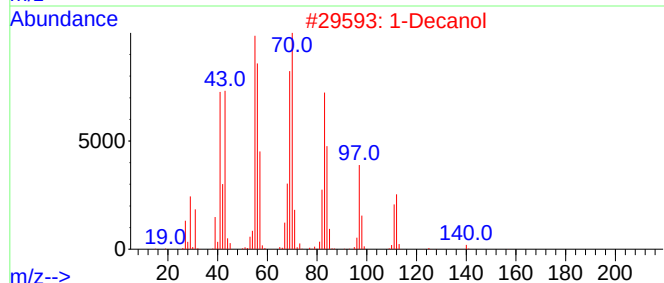
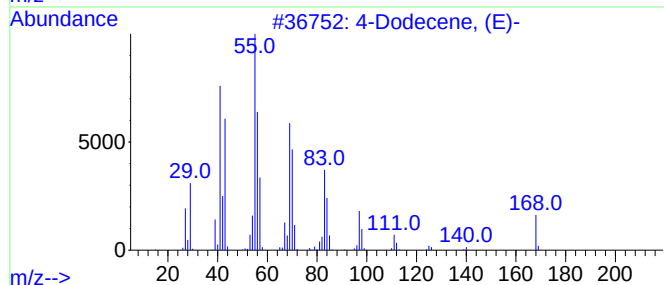
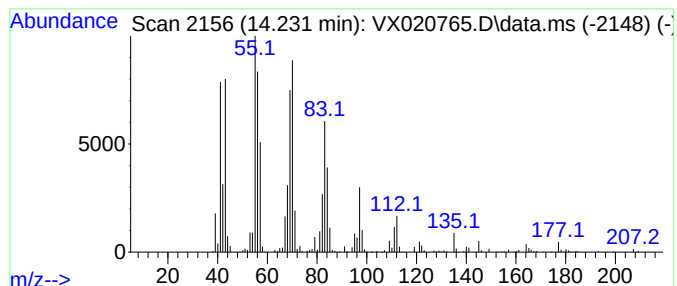
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TIC Integration Parameters: LSCINT.P

Peak Number 4 4-Dodecene, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.231	12.12 ug/l	410738	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Dodecene, (E)-	168	C12H24	007206-15-7	94
2			1-Decanol	158	C10H22O	000112-30-1	91
3			1-Dodecanol	186	C12H26O	000112-53-8	91
4			Cyclooctane	112	C8H16	000292-64-8	87
5			Cyclopropane, nonyl-	168	C12H24	074663-85-7	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012621\
Data File : VX020765.D
Acq On : 26 Jan 2021 19:48
Operator : JC/SP
Sample : M1217-16 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1710

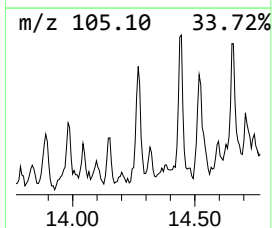
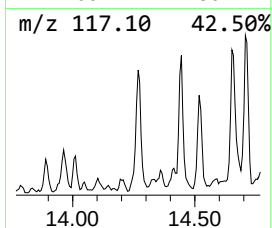
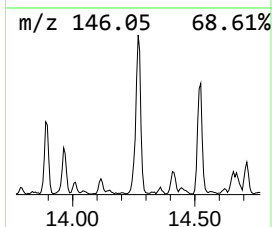
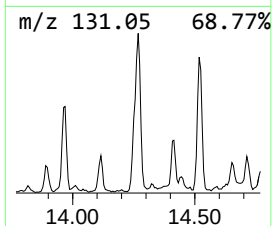
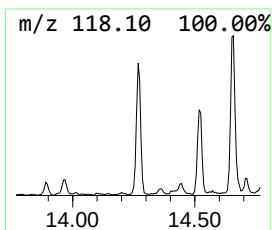
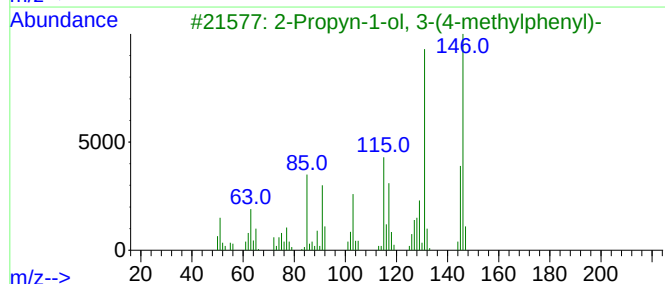
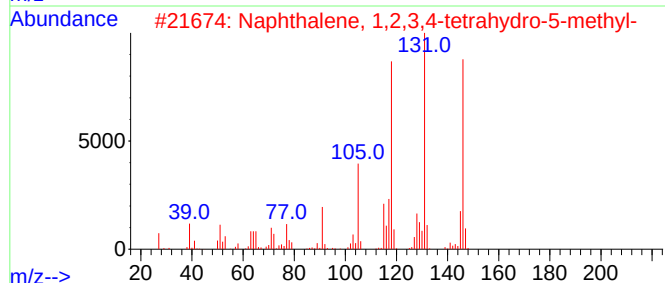
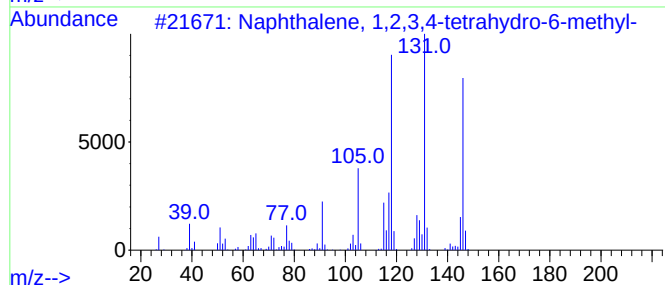
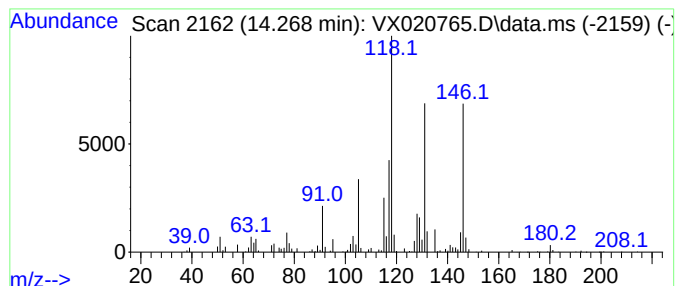
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Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.268	10.81 ug/l	366649	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	91
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	87
3			2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	50
4			1(2H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000529-34-0	49
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012621\
Data File : VX020765.D
Acq On : 26 Jan 2021 19:48
Operator : JC/SP
Sample : M1217-16 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1710

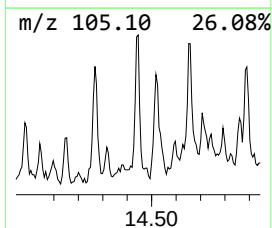
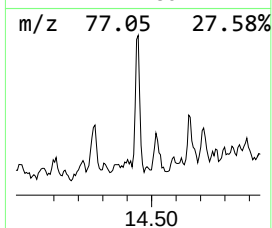
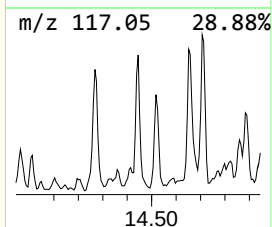
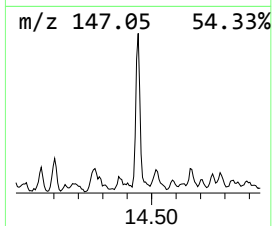
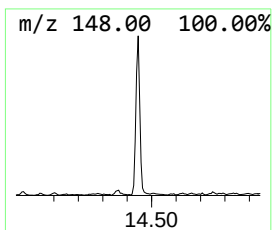
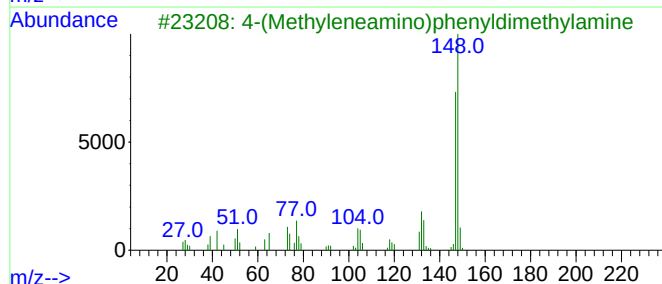
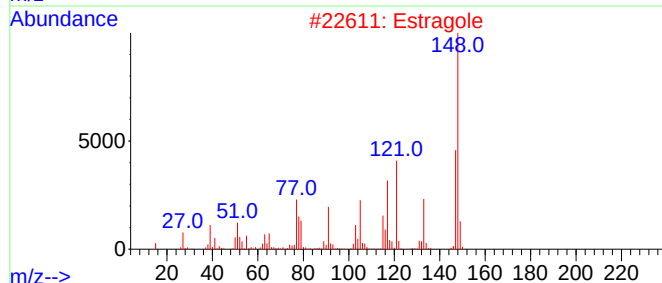
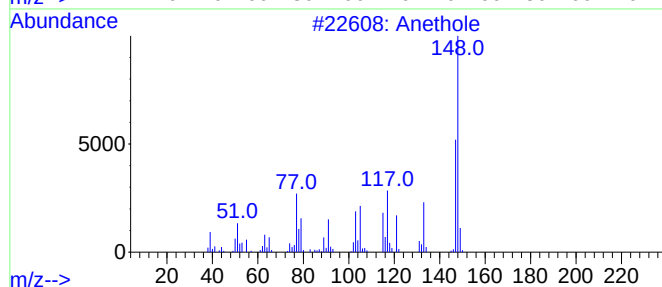
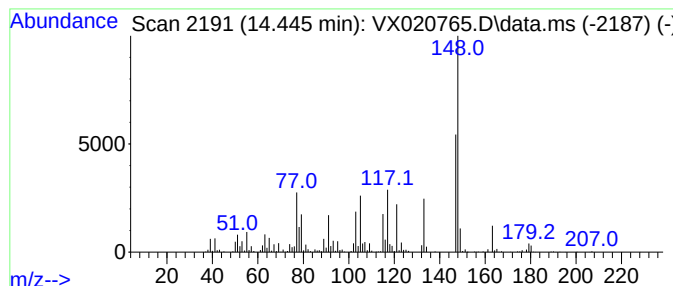
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Quant Title : SW846 8260

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

Peak Number 6 Anethole Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.445	9.04 ug/l	306551	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anethole			148	C10H12O	000104-46-1	93
2	Estragole			148	C10H12O	000140-67-0	87
3	4-(Methyleneamino)phenyldimethyl...			148	C9H12N2	147354-14-1	52
4	2-(1-Propenyl)-6-methylphenol			148	C10H12O	1000306-52-9	50
5	2-Hydroxymethylbenzimidazole			148	C8H8N2O	022128-99-0	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012621\
Data File : VX020765.D
Acq On : 26 Jan 2021 19:48
Operator : JC/SP
Sample : M1217-16 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

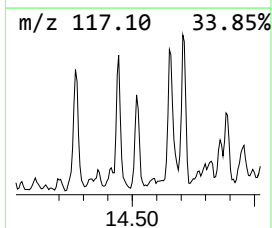
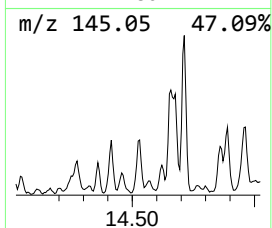
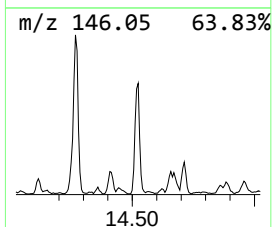
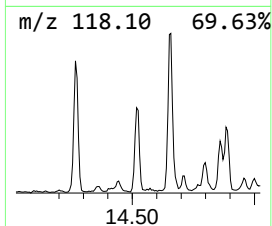
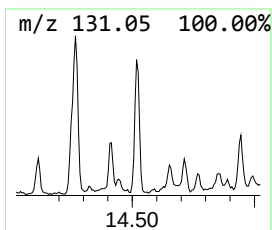
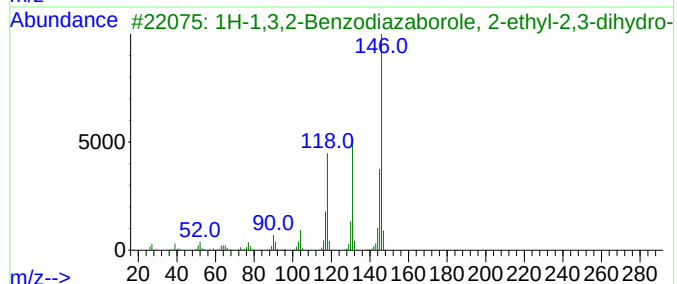
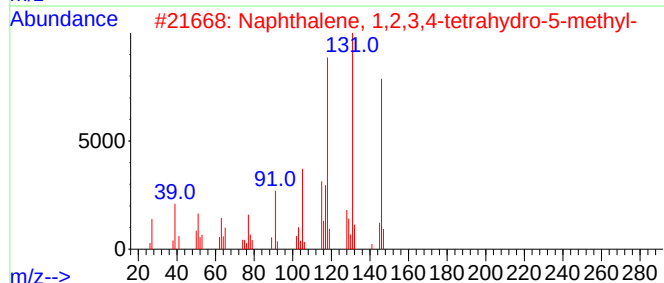
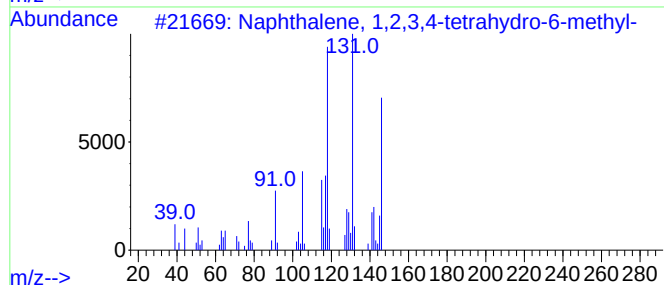
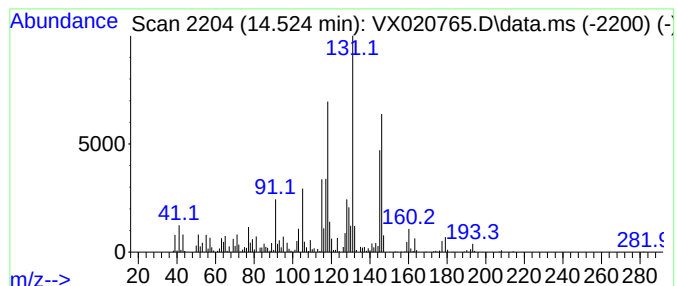
Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X012521W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.524	10.24 ug/l	347213	1,4-Dichlorobenzene-d4	12.061	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	96
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
3	1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	70
4	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	64
5	1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012621\
Data File : VX020765.D
Acq On : 26 Jan 2021 19:48
Operator : JC/SP
Sample : M1217-16 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1710

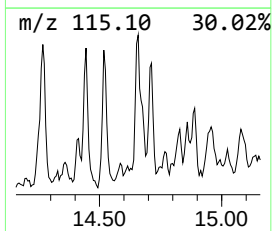
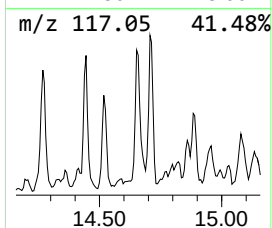
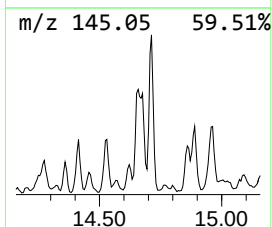
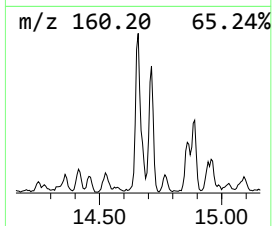
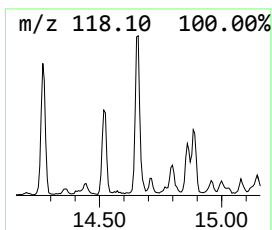
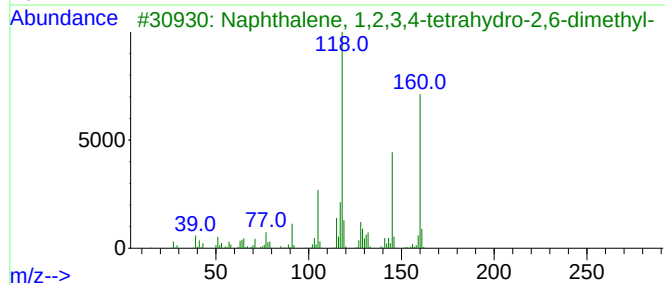
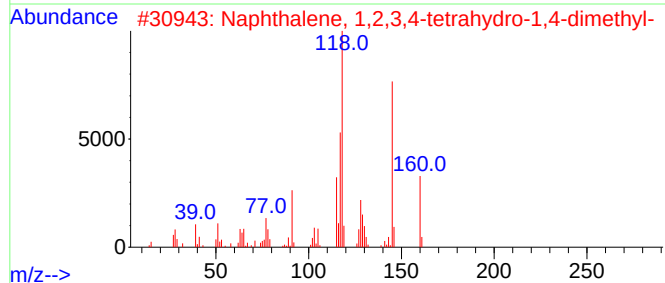
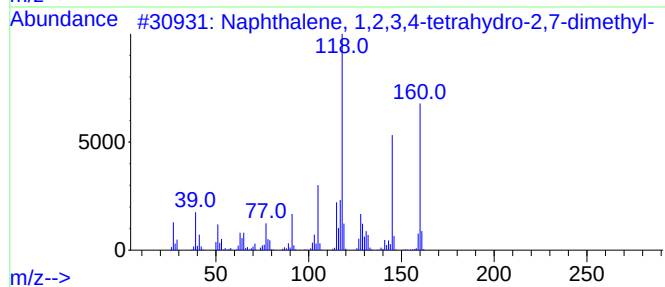
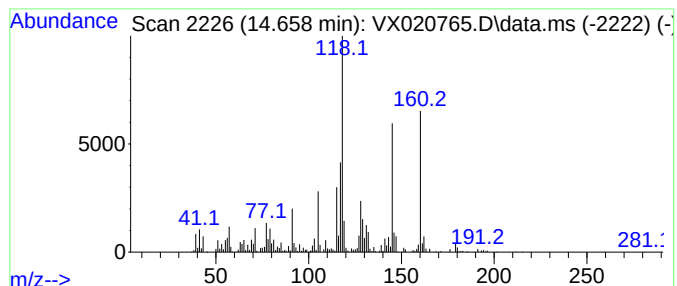
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.658	12.84 ug/l	435329	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	90
2			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	90
3			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	83
4			Benzene, (3-methyl-1-methylenebu...	160	C12H16	038212-14-5	64
5			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012621\
Data File : VX020765.D
Acq On : 26 Jan 2021 19:48
Operator : JC/SP
Sample : M1217-16 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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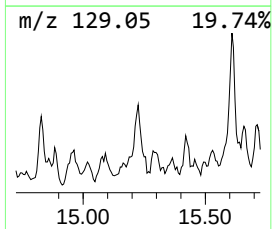
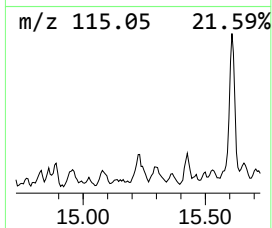
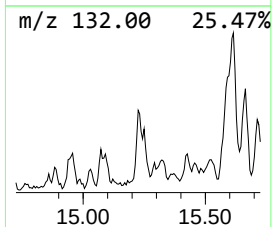
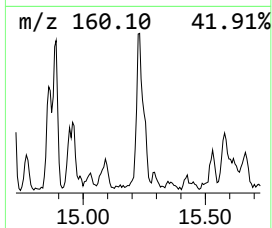
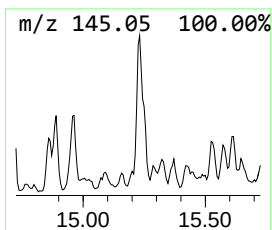
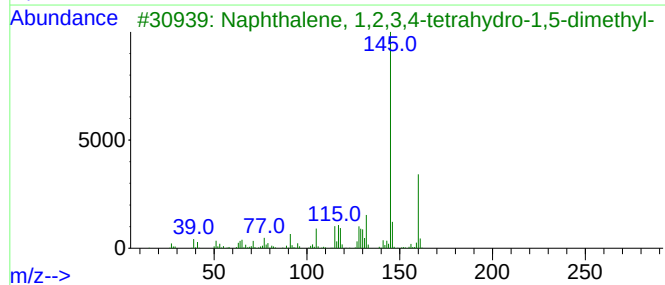
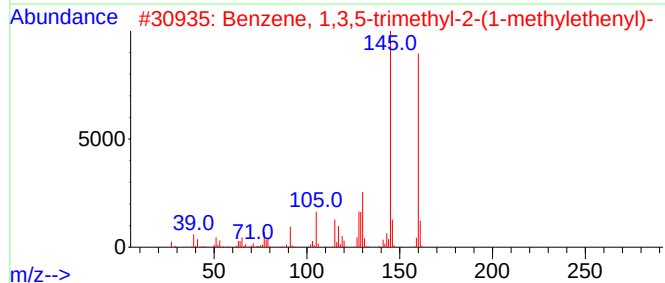
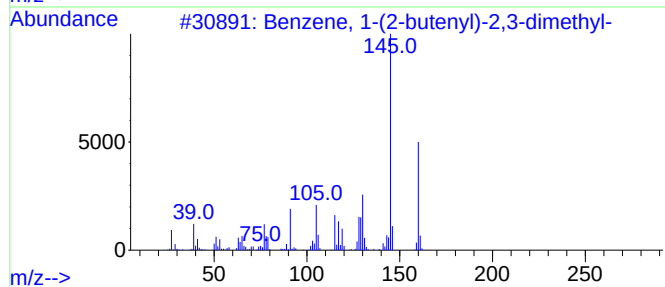
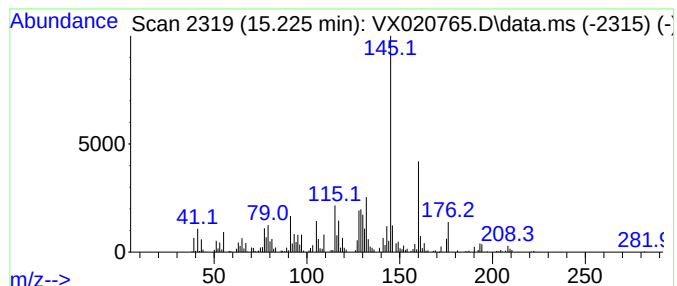
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-(2-butenyl)-2,3-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.225	8.03 ug/l	272121	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	91
2			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	90
3			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	87
4			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	76
5			Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012621\
Data File : VX020765.D
Acq On : 26 Jan 2021 19:48
Operator : JC/SP
Sample : M1217-16 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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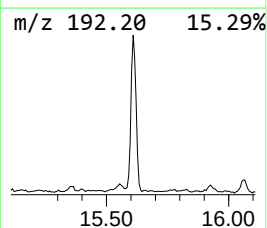
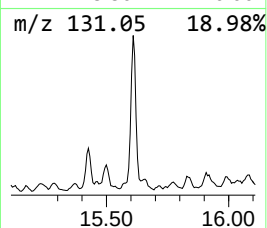
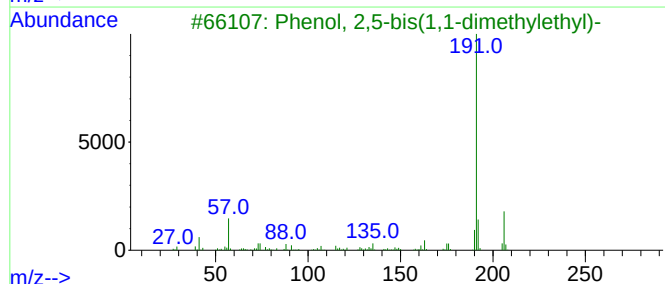
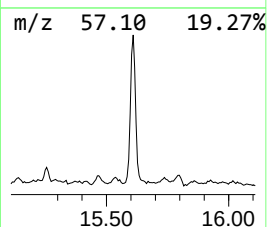
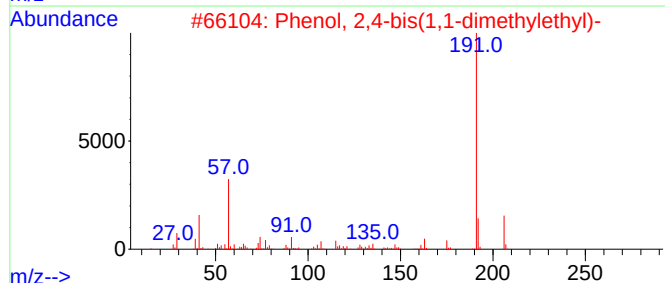
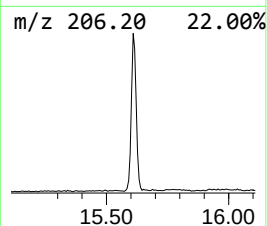
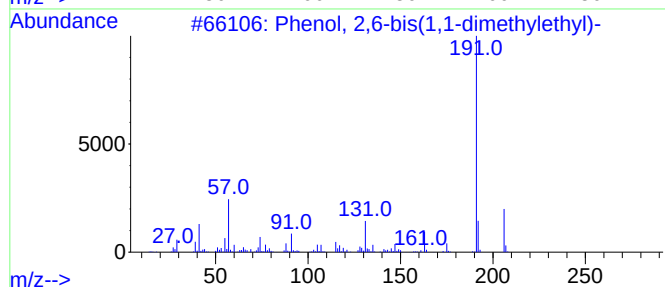
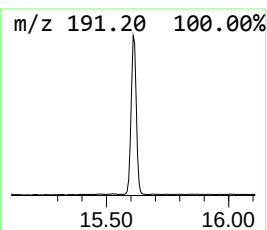
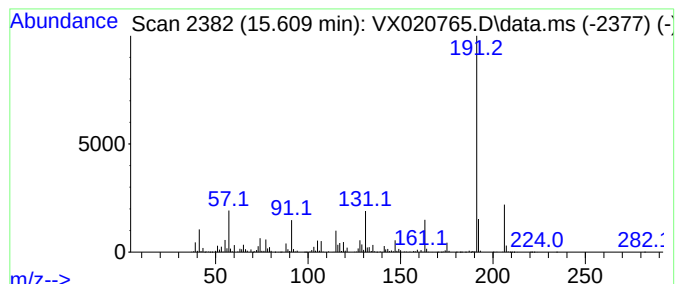
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Quant Title : SW846 8260

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

Peak Number 10 Phenol, 2,6-bis(1,1-dimethy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.609	52.92 ug/l	1794200	1,4-Dichlorobenzene-d4	12.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	96
2			Phenol, 2,4-bis(1,1-dimethylethyl)-	206	C14H22O	000096-76-4	87
3			Phenol, 2,5-bis(1,1-dimethylethyl)-	206	C14H22O	005875-45-6	81
4			Phenol, 3,5-bis(1,1-dimethylethyl)-	206	C14H22O	001138-52-9	81
5			Ethyl 4-t-butylbenzoate	206	C13H18O2	005406-57-5	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012621\
Data File : VX020765.D
Acq On : 26 Jan 2021 19:48
Operator : JC/SP
Sample : M1217-16 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X012521W.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
Eucalyptol	12.140	44.0	ug/l	1492060	4	12.061	1695140	50.0
Naphthalene, 1,...	13.457	35.1	ug/l	1189150	4	12.061	1695140	50.0
Naphthalene, 1,...	13.890	9.0	ug/l	305434	4	12.061	1695140	50.0
4-Dodecene, (E)-	14.231	12.1	ug/l	410738	4	12.061	1695140	50.0
Naphthalene, 1,...	14.268	10.8	ug/l	366649	4	12.061	1695140	50.0
Anethole	14.445	9.0	ug/l	306551	4	12.061	1695140	50.0
Naphthalene, 1,...	14.524	10.2	ug/l	347213	4	12.061	1695140	50.0
Naphthalene, 1,...	14.658	12.8	ug/l	435329	4	12.061	1695140	50.0
Benzene, 1-(2-b...	15.225	8.0	ug/l	272121	4	12.061	1695140	50.0
Phenol, 2,6-bis...	15.609	52.9	ug/l	1794200	4	12.061	1695140	50.0