

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW062121\
 Data File : VW019252.D
 Acq On : 21 Jun 2021 16:14
 Operator : SY/VA
 Sample : M2777-03
 Misc : 3.71G/5ML/MSVOA_W/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 X-3

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_W\Method\82W062121S.M
 Title : SW846 8260

Signal : TIC: VW019252.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.921	484	495	513	rBV3	20006	79811	1.89%	0.130%
2	7.141	850	859	871	rBV	28031	80991	1.91%	0.132%
3	7.890	972	982	986	rBV	250033	583894	13.80%	0.954%
4	7.951	986	992	1003	rVB	436054	978261	23.13%	1.599%
5	8.311	1034	1051	1062	rBV	268037	622324	14.71%	1.017%
6	8.842	1129	1138	1151	rBV	679289	1352408	31.97%	2.210%
7	10.323	1374	1381	1396	rBV	1068271	1918022	45.35%	3.134%
8	11.061	1489	1502	1506	rBV9	32598	120002	2.84%	0.196%
9	11.231	1526	1530	1537	rVB3	31329	61801	1.46%	0.101%
10	11.341	1543	1548	1552	rBV4	24084	44881	1.06%	0.073%
11	11.512	1572	1576	1581	rBV2	33636	60314	1.43%	0.099%
12	11.634	1589	1596	1607	rVB	917415	1594194	37.69%	2.605%
13	11.841	1627	1630	1633	rVV	54611	72032	1.70%	0.118%
14	12.140	1673	1679	1690	rBV3	72194	195495	4.62%	0.319%
15	12.249	1692	1697	1702	rVB	82605	122122	2.89%	0.200%
16	12.365	1708	1716	1722	rBV5	77271	199055	4.71%	0.325%
17	12.615	1752	1757	1767	rVB	743052	1278948	30.24%	2.090%
18	12.780	1780	1784	1791	rVB4	80978	148907	3.52%	0.243%
19	12.859	1791	1797	1801	rBV4	51636	124398	2.94%	0.203%
20	12.932	1801	1809	1816	rBV3	328903	718998	17.00%	1.175%
21	12.993	1816	1819	1822	rVB2	39667	49216	1.16%	0.080%
22	13.079	1823	1833	1834	rBV3	132089	319449	7.55%	0.522%
23	13.164	1843	1847	1855	rVB2	185049	328219	7.76%	0.536%
24	13.298	1865	1869	1873	rBV4	66244	129667	3.07%	0.212%
25	13.377	1880	1882	1886	rVB2	64092	76396	1.81%	0.125%
26	13.438	1886	1892	1897	rBV2	276687	511194	12.09%	0.835%
27	13.487	1897	1900	1903	rVV2	103134	145201	3.43%	0.237%
28	13.554	1906	1911	1920	rVB2	859821	1560277	36.89%	2.550%
29	13.621	1920	1922	1928	rVB3	81463	103535	2.45%	0.169%
30	13.719	1929	1938	1939	rBV6	111602	248954	5.89%	0.407%
31	13.749	1939	1943	1946	rVV	196533	289753	6.85%	0.474%
32	13.786	1946	1949	1954	rVB2	316684	456068	10.78%	0.745%
33	13.871	1958	1963	1969	rBV3	647612	1114006	26.34%	1.820%
34	13.950	1972	1976	1983	rVB	258081	407430	9.63%	0.666%
35	14.097	1995	2000	2004	rVB	404011	598708	14.15%	0.978%
36	14.133	2004	2006	2011	rVB2	84466	94212	2.23%	0.154%

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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_W\Method\82W062121S.M
 Title : SW846 8260

37	14.200	2011	2017	2022	rBV	721040	1288329	30.46%	2.105%
38	14.255	2022	2026	2028	rBV2	130691	198983	4.70%	0.325%
39	14.341	2034	2040	2043	rBV2	258678	500338	11.83%	0.818%
40	14.383	2043	2047	2050	rVV2	362165	587235	13.88%	0.960%
41	14.420	2050	2053	2056	rVV2	278981	427707	10.11%	0.699%
42	14.456	2056	2059	2062	rVB	441954	541295	12.80%	0.885%
43	14.499	2062	2066	2069	rBV	461070	565947	13.38%	0.925%
44	14.542	2069	2073	2080	rVB3	384933	727506	17.20%	1.189%
45	14.639	2080	2089	2093	rBV5	209743	514849	12.17%	0.841%
46	14.682	2093	2096	2099	rBV3	187132	278334	6.58%	0.455%
47	14.725	2099	2103	2108	rVB2	912556	1402265	33.15%	2.292%
48	14.804	2108	2116	2121	rBV3	1064258	3004683	71.04%	4.910%
49	14.853	2121	2124	2129	rVB2	343745	611311	14.45%	0.999%
50	14.956	2137	2141	2148	rVB2	639445	1055526	24.95%	1.725%
51	15.023	2149	2152	2153	rBV3	113683	126385	2.99%	0.207%
52	15.060	2153	2158	2163	rBV2	706382	1411572	33.37%	2.307%
53	15.176	2170	2177	2184	rBV2	791417	1902040	44.97%	3.108%
54	15.243	2185	2188	2196	rVB6	282465	539541	12.76%	0.882%
55	15.322	2196	2201	2206	rBV5	402571	814950	19.27%	1.332%
56	15.420	2211	2217	2221	rBV	1007736	1761622	41.65%	2.879%
57	15.499	2221	2230	2235	rVV3	768760	2411994	57.02%	3.942%
58	15.548	2235	2238	2247	rVB2	808957	1461684	34.56%	2.389%
59	15.651	2248	2255	2261	rBV3	755184	1674229	39.58%	2.736%
60	15.724	2264	2267	2272	rVB3	166476	240688	5.69%	0.393%
61	15.822	2275	2283	2285	rVV4	576872	1316088	31.12%	2.151%
62	15.859	2285	2289	2299	rVV	1547091	2995205	70.81%	4.895%
63	15.950	2299	2304	2309	rVV4	284135	581807	13.76%	0.951%
64	16.023	2309	2316	2330	rVB4	618959	2177891	51.49%	3.559%
65	16.176	2331	2341	2353	rBV	1517864	3797165	89.77%	6.205%
66	16.279	2355	2358	2363	rVB2	179967	289886	6.85%	0.474%
67	16.371	2363	2373	2379	rBV2	1836933	4229748	100.00%	6.912%
68	16.438	2379	2384	2388	rVV	565229	1135104	26.84%	1.855%
69	16.486	2388	2392	2397	rVB2	673017	1218355	28.80%	1.991%
70	16.651	2409	2419	2423	rBV4	562444	1767308	41.78%	2.888%
71	16.779	2434	2440	2441	rBV	528749	846267	20.01%	1.383%

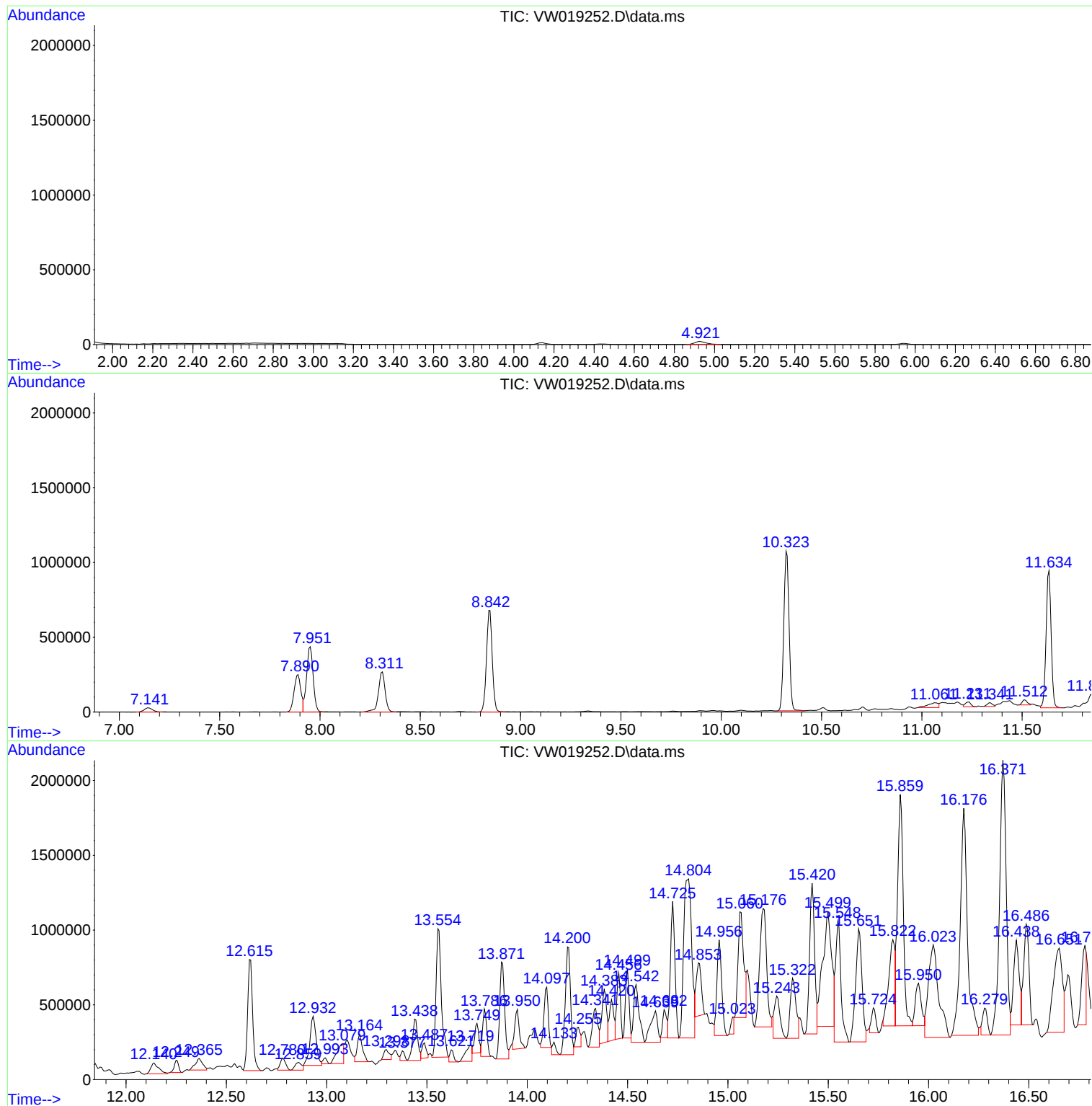
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ClientSampled :
X-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\82W062121S.M
Quant Title : SW846 8260

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



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Misc : 3.71G/5ML/MSVOA_W/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampled :
X-3

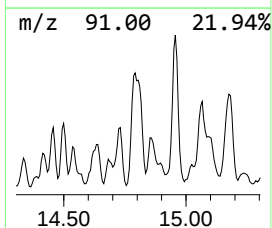
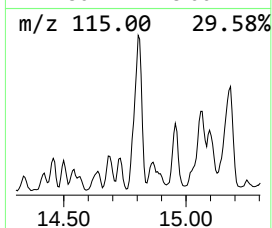
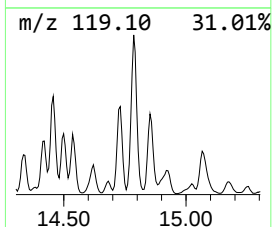
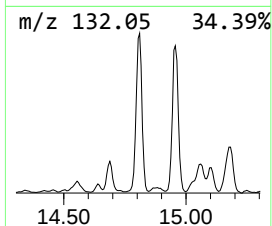
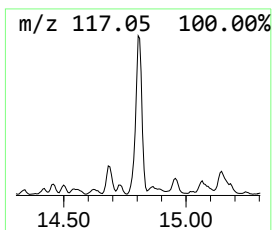
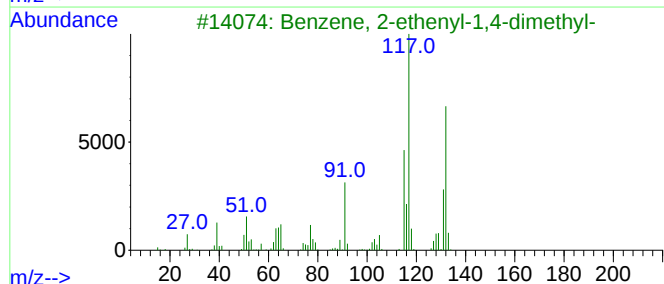
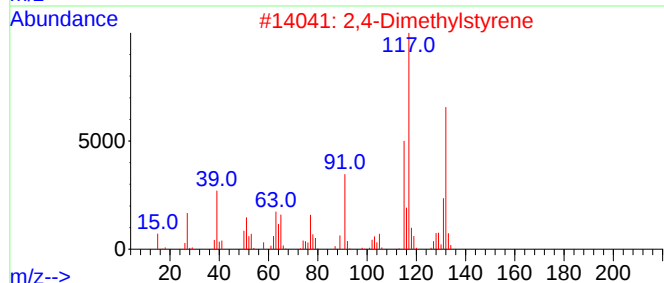
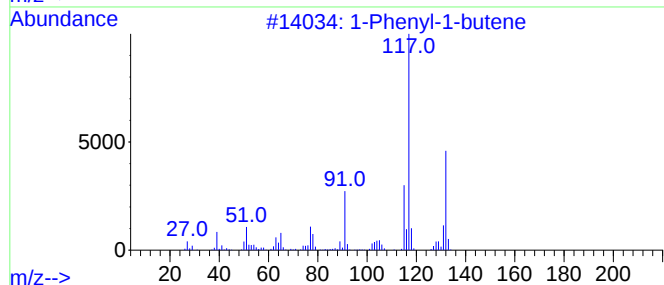
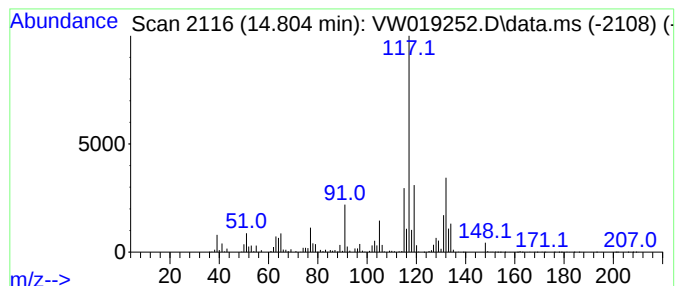
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\82W062121S.M
Quant Title : SW846 8260

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

Peak Number 1 1-Phenyl-1-butene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.804	96.29 ug/l	3004680	1,4-Dichlorobenzene-d4	13.560

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Phenyl-1-butene	132	C10H12	000824-90-8	95
2		2,4-Dimethylstyrene	132	C10H12	002234-20-0	91
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
4		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	87
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	76



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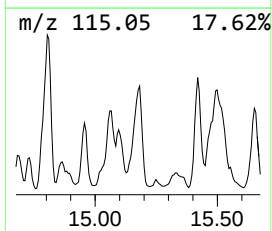
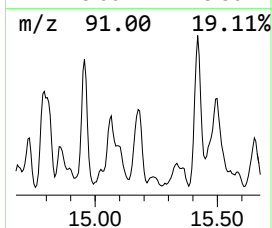
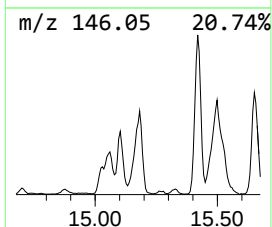
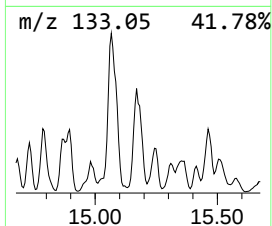
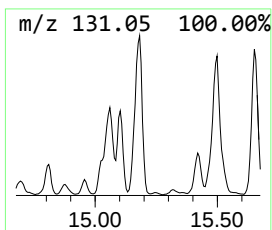
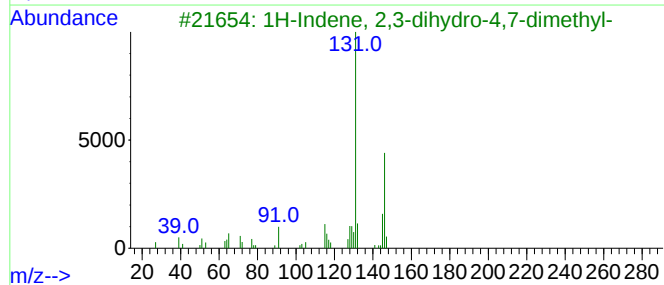
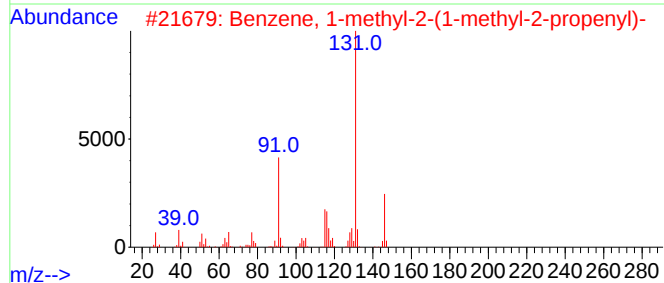
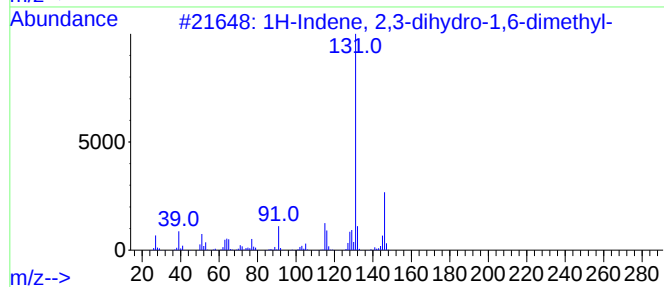
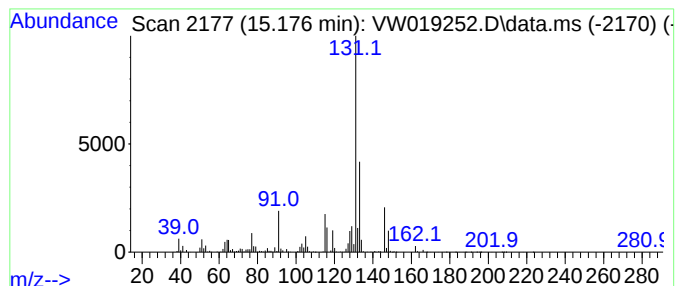
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\82W062121S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.176	60.95 ug/l	1902040	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91
2			Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	70
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64
4			1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	64
5			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	64



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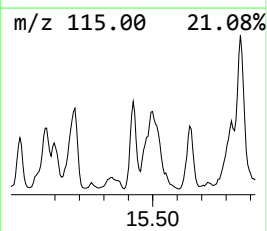
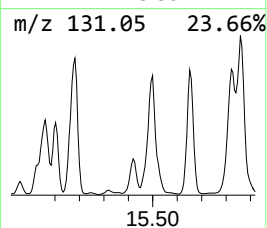
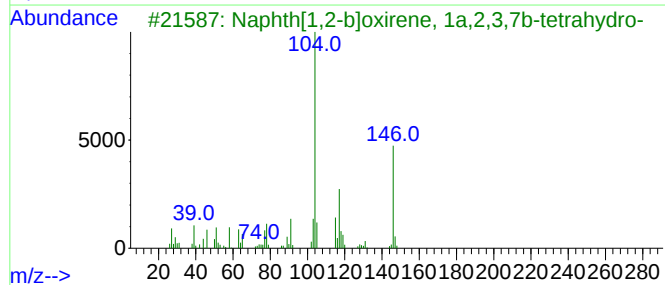
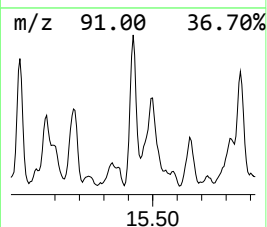
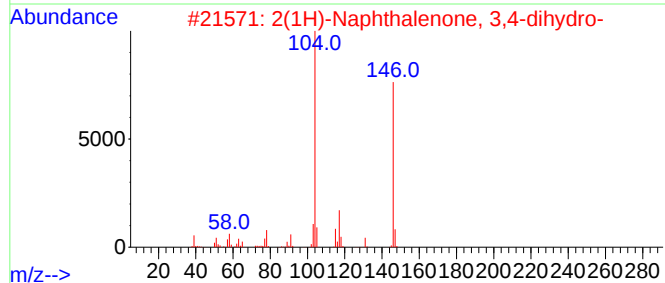
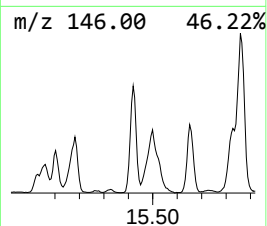
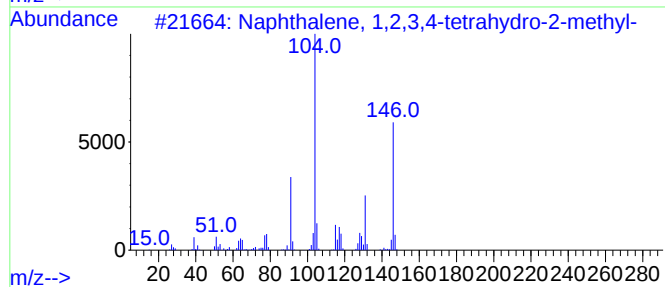
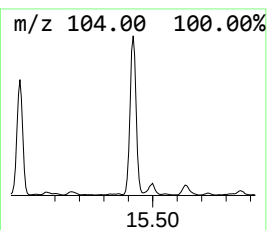
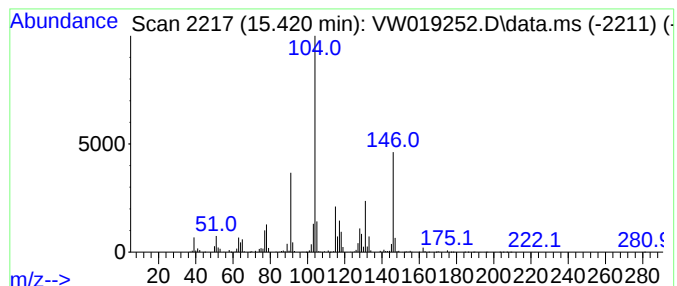
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.420	56.45 ug/l	1761620	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	96
2			2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	68
3			Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	62
4			2-Azetidinone, 4-phenyl-	147	C9H9NO	005661-55-2	43
5			Benzene, 3-cyclohexen-1-yl-	158	C12H14	004994-16-5	38



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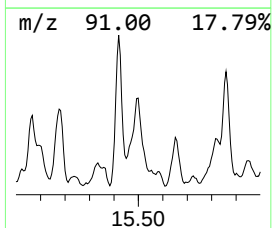
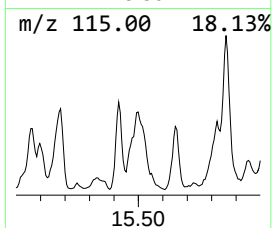
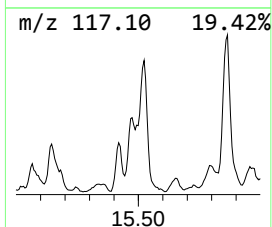
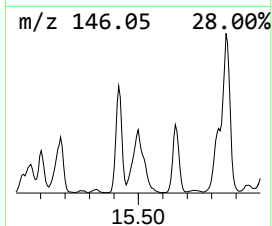
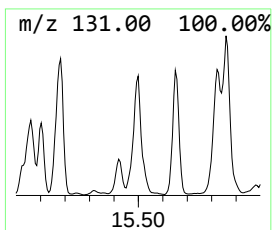
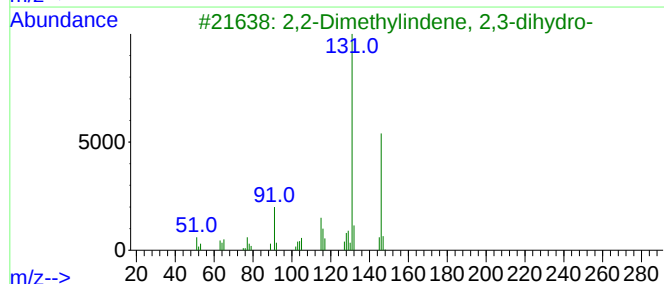
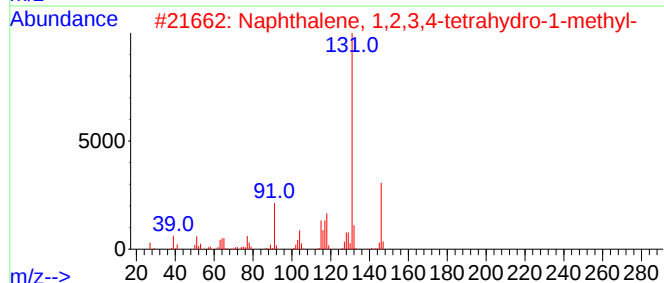
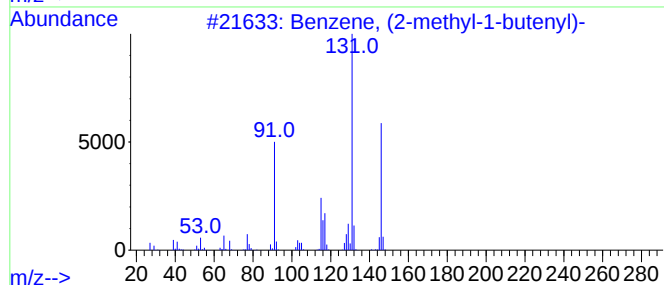
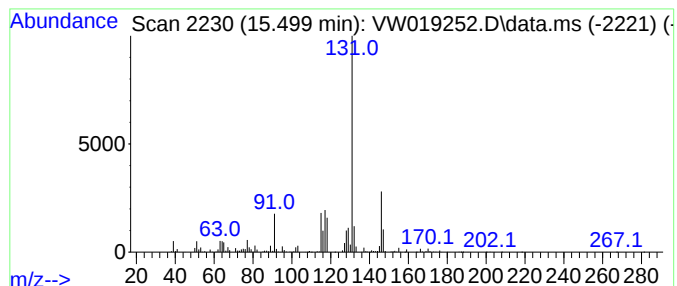
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\82W062121S.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, (2-methyl-1-butenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.499	77.29 ug/l	2411990	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	93
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	90
3			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	81
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	74
5			1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW062121\
Data File : VW019252.D
Acq On : 21 Jun 2021 16:14
Operator : SY/VA
Sample : M2777-03
Misc : 3.71G/5ML/MSVOA_W/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
X-3

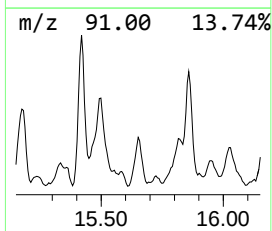
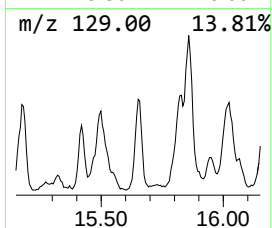
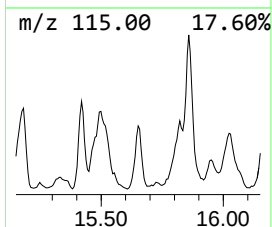
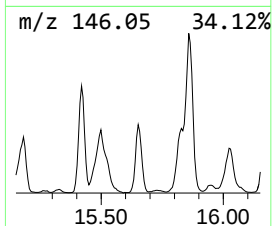
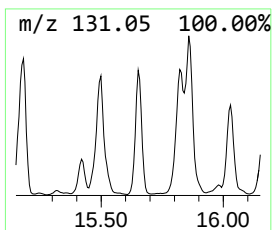
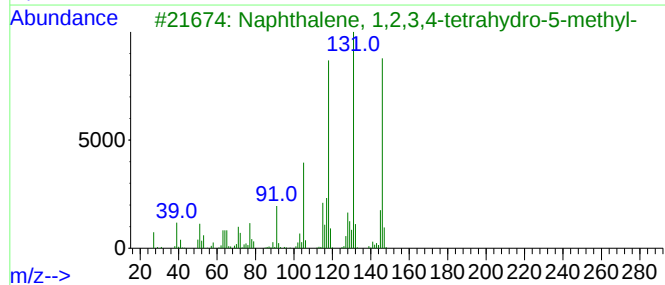
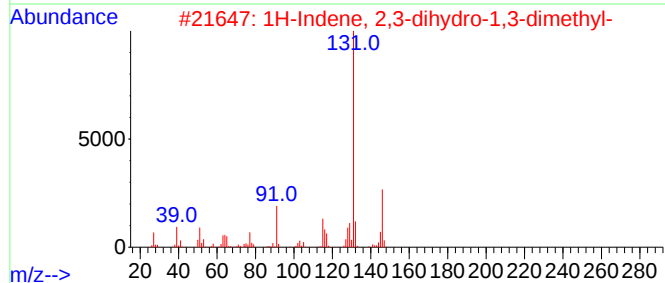
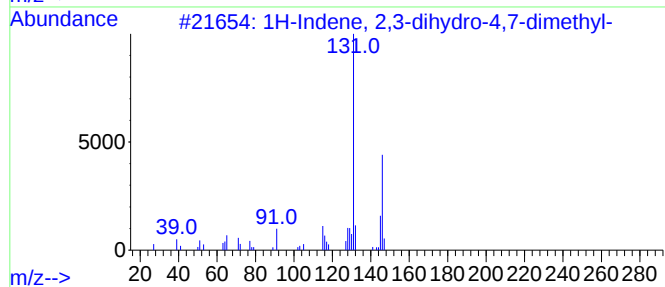
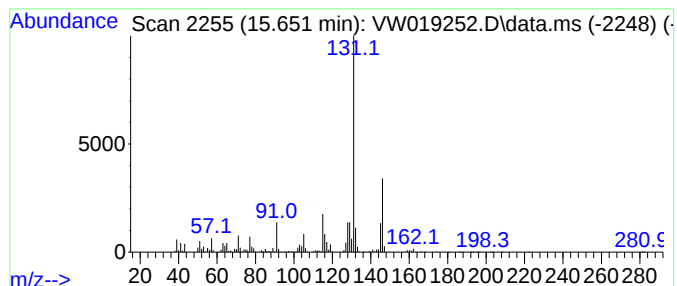
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Quant Title : SW846 8260

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

Peak Number 5 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.651	53.65 ug/l	1674230	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
2			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	91
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
5			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW062121\
Data File : VW019252.D
Acq On : 21 Jun 2021 16:14
Operator : SY/VA
Sample : M2777-03
Misc : 3.71G/5ML/MSVOA_W/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampled :
X-3

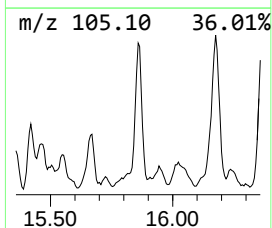
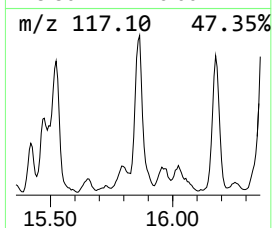
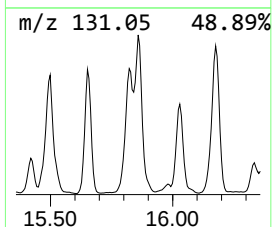
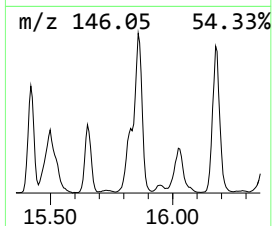
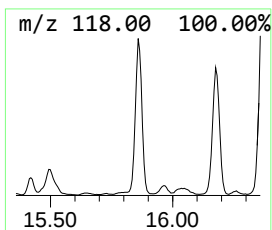
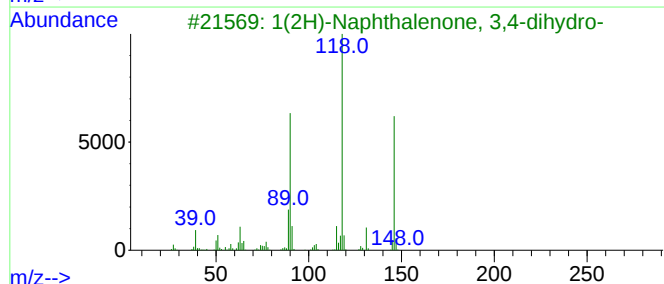
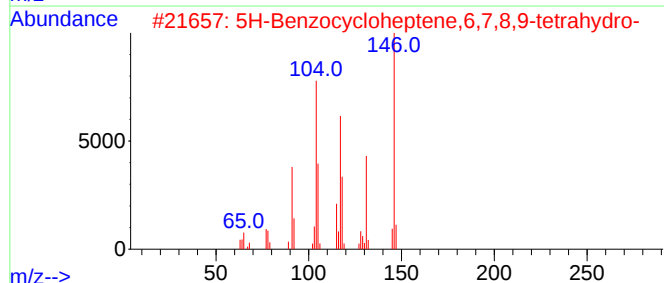
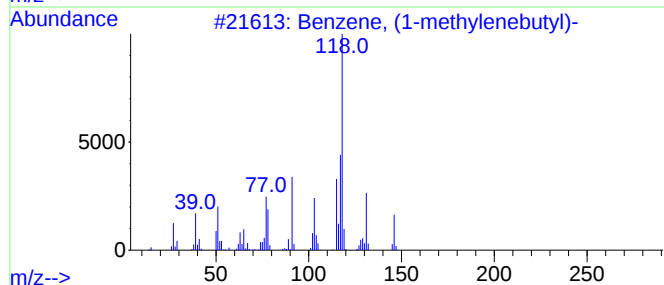
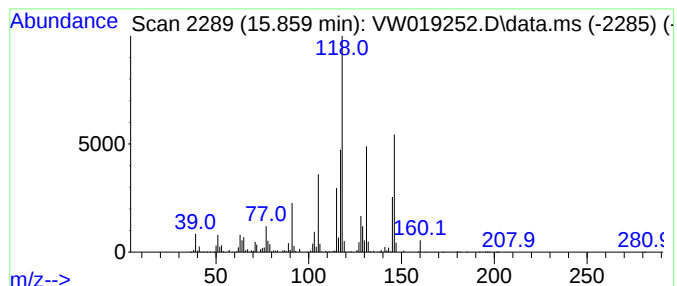
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, (1-methylenebutyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.858	95.98 ug/l	2995210	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	80
2			5H-Benzocycloheptene,6,7,8,9-tet...	146	C11H14	001075-16-7	47
3			1(2H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000529-34-0	38
4			trans-Tetraline-2,3-diol	164	C10H12O2	041597-55-1	32
5			Cinnamaldehyde, .beta.-methyl-	146	C10H10O	001196-67-4	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW062121\
Data File : VW019252.D
Acq On : 21 Jun 2021 16:14
Operator : SY/VA
Sample : M2777-03
Misc : 3.71G/5ML/MSVOA_W/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
X-3

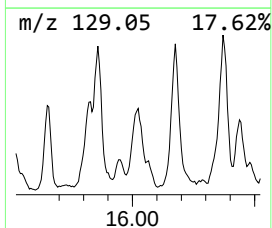
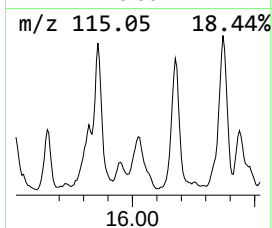
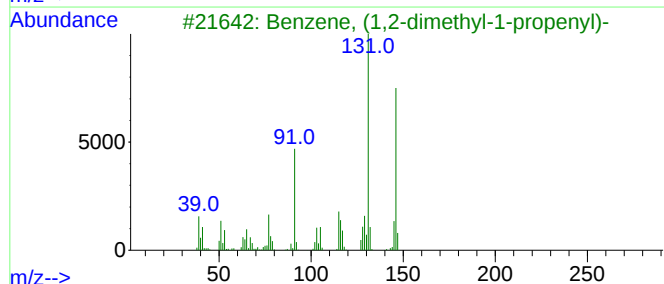
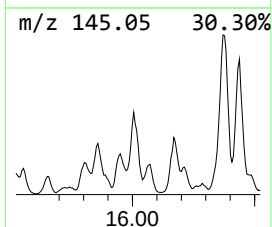
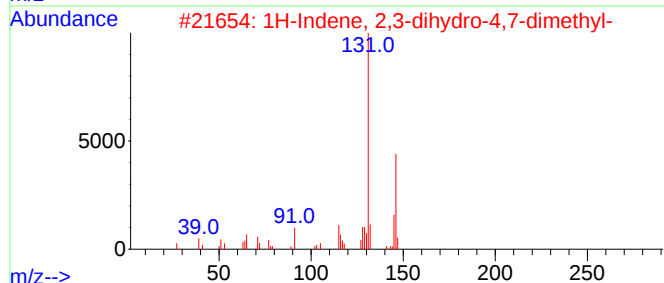
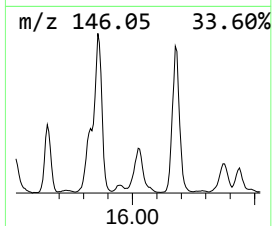
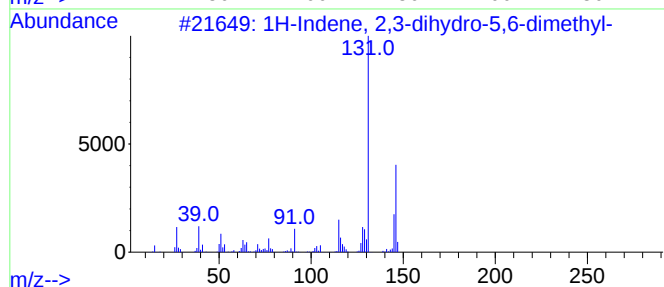
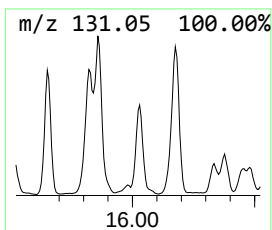
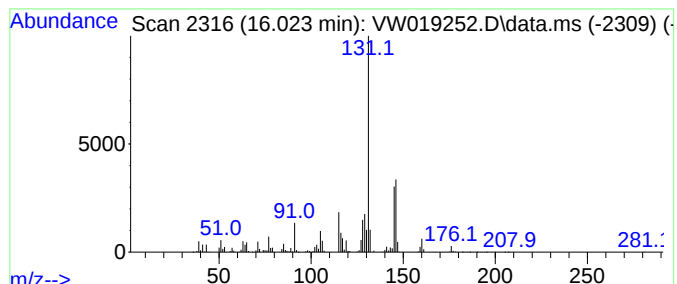
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Quant Title : SW846 8260

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

Peak Number 7 1H-Indene, 2,3-dihydro-5,6-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.023	69.79 ug/l	2177890	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	93
2			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
3			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	87
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	87
5			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW062121\
Data File : VW019252.D
Acq On : 21 Jun 2021 16:14
Operator : SY/VA
Sample : M2777-03
Misc : 3.71G/5ML/MSVOA_W/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
X-3

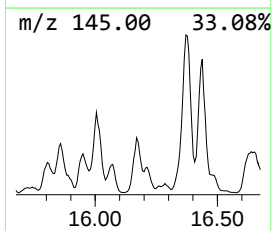
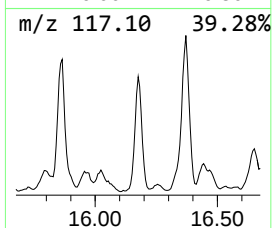
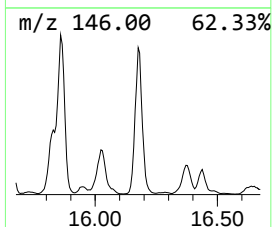
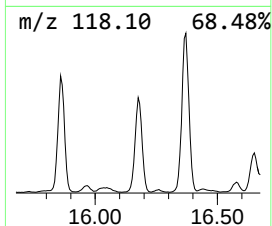
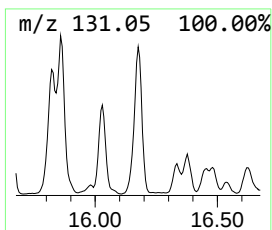
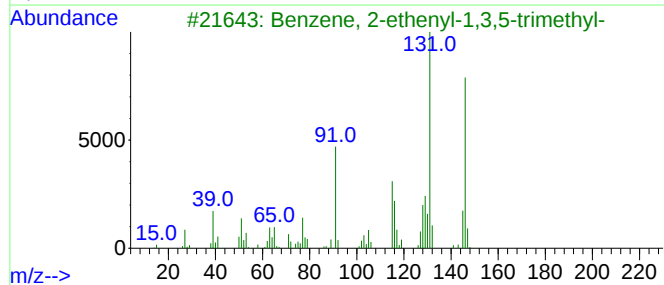
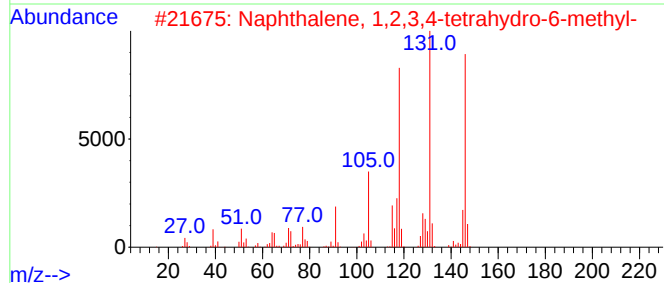
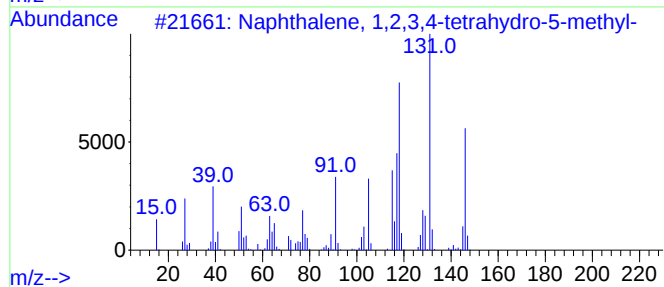
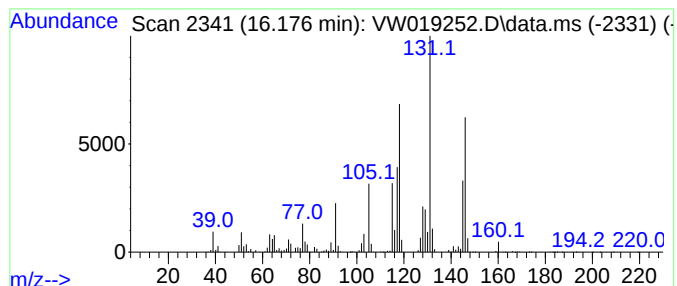
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\82W062121S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.176	121.68 ug/l	3797170	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	93
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	81
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	78
4			1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	68
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW062121\
Data File : VW019252.D
Acq On : 21 Jun 2021 16:14
Operator : SY/VA
Sample : M2777-03
Misc : 3.71G/5ML/MSVOA_W/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
X-3

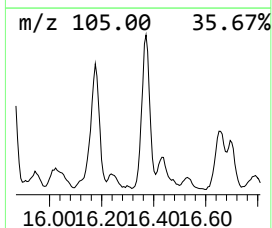
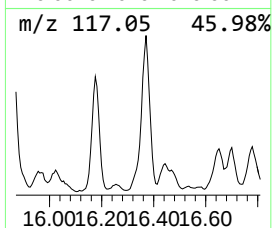
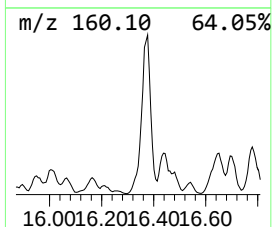
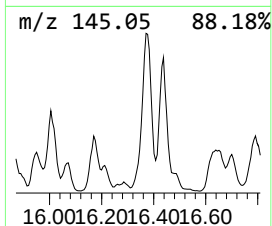
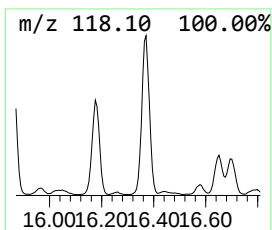
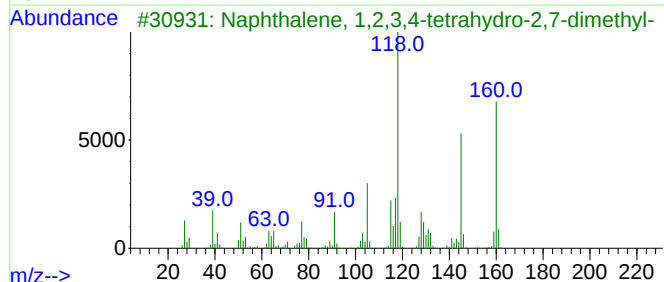
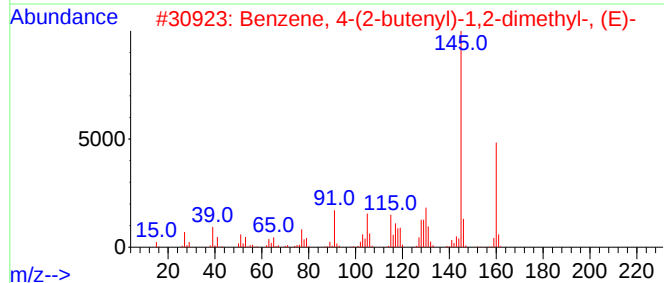
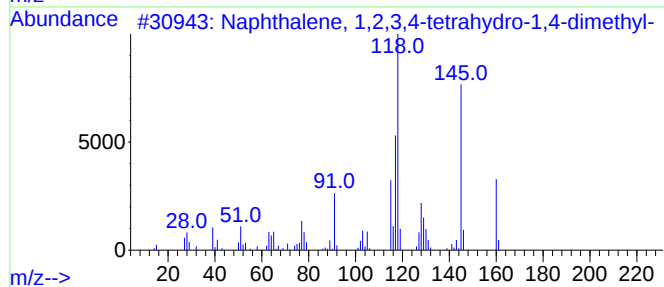
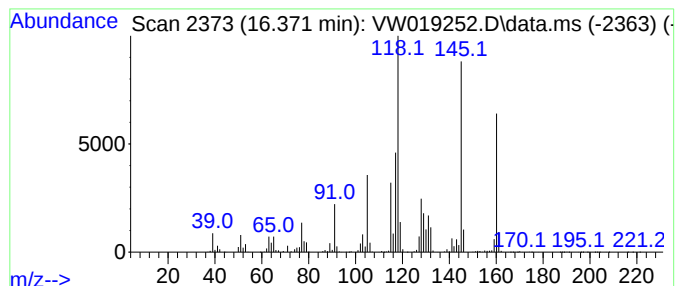
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\82W062121S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.371	135.54 ug/l	4229750	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	95
2			Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	83
3			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	70
4			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	70
5			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW062121\
Data File : VW019252.D
Acq On : 21 Jun 2021 16:14
Operator : SY/VA
Sample : M2777-03
Misc : 3.71G/5ML/MSVOA_W/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
X-3

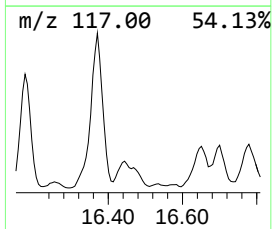
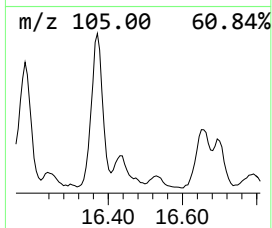
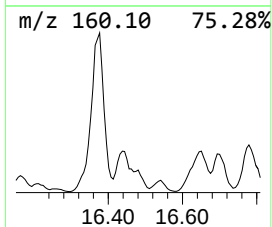
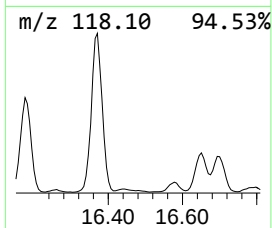
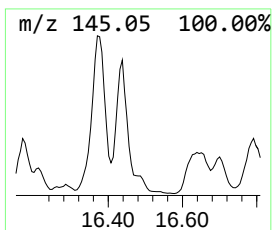
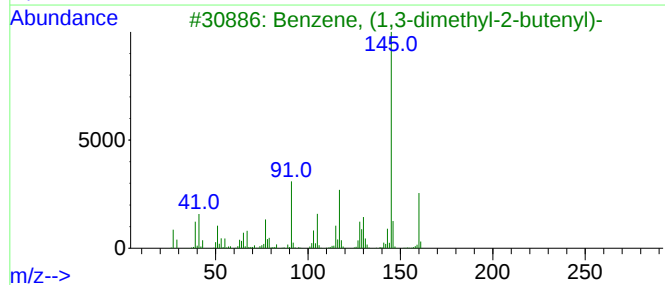
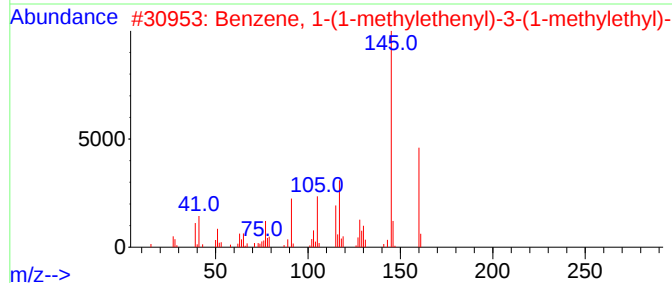
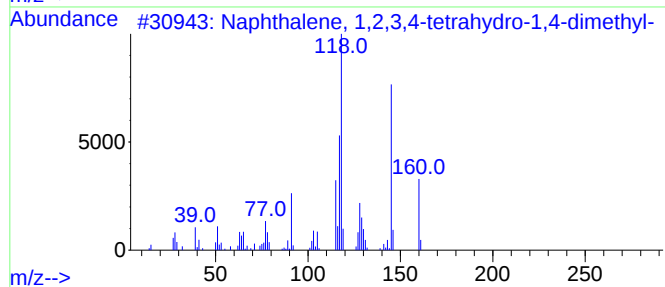
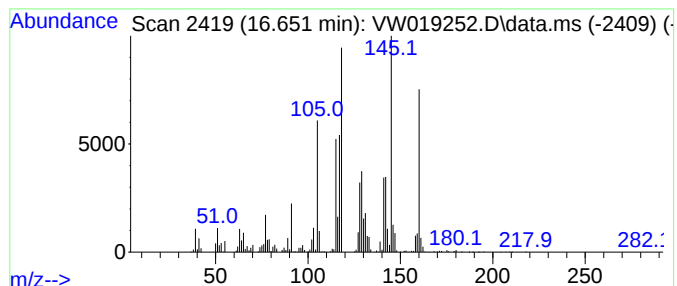
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\82W062121S.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-(1-methylethenyl)... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.651	56.63 ug/l	1767310	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	81
2			Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	64
3			Benzene, (1,3-dimethyl-2-butenyl)-	160	C12H16	050704-01-3	53
4			1H-Indene, 2,3-dihydro-1,5,7-tri...	160	C12H16	054340-88-4	49
5			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW062121\
 Data File : VW019252.D
 Acq On : 21 Jun 2021 16:14
 Operator : SY/VA
 Sample : M2777-03
 Misc : 3.71G/5ML/MSVOA_W/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 X-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\82W062121S.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
1-Phenyl-1-butene	14.804	96.3	ug/l	3004680	4	13.560	1560280	50.0
1H-Indene, 2,3-...	15.176	61.0	ug/l	1902040	4	13.560	1560280	50.0
Naphthalene, 1,...	15.420	56.5	ug/l	1761620	4	13.560	1560280	50.0
Benzene, (2-met...	15.499	77.3	ug/l	2411990	4	13.560	1560280	50.0
1H-Indene, 2,3-...	15.651	53.6	ug/l	1674230	4	13.560	1560280	50.0
Benzene, (1-met...	15.858	96.0	ug/l	2995210	4	13.560	1560280	50.0
1H-Indene, 2,3-...	16.023	69.8	ug/l	2177890	4	13.560	1560280	50.0
Naphthalene, 1,...	16.176	121.7	ug/l	3797170	4	13.560	1560280	50.0
Naphthalene, 1,...	16.371	135.5	ug/l	4229750	4	13.560	1560280	50.0
Benzene, 1-(1-m...	16.651	56.6	ug/l	1767310	4	13.560	1560280	50.0