

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD061621\
 Data File : VD069461.D
 Acq On : 16 Jun 2021 12:32
 Operator : VA/SY
 Sample : M2709-01
 Misc : 5.00G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 0228-COMPOSITE-E

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_D\Method\82D061121S.M

Title : SW846 8260

Signal : TIC: VD069461.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.156	372	380	388	rBV3	21234	63050	1.46%	0.147%
2	4.967	507	518	528	rBV4	29918	96451	2.24%	0.224%
3	7.202	886	898	906	rBV3	22164	68916	1.60%	0.160%
4	7.914	1008	1019	1024	rBV	245312	589972	13.70%	1.372%
5	7.979	1024	1030	1039	rVV	357813	813642	18.90%	1.892%
6	8.061	1039	1044	1054	rVB4	26528	65016	1.51%	0.151%
7	8.332	1077	1090	1102	rVB2	208569	504525	11.72%	1.173%
8	8.455	1102	1111	1114	rVV6	17580	44120	1.02%	0.103%
9	8.508	1115	1120	1129	rVV4	28971	94545	2.20%	0.220%
10	8.602	1129	1136	1144	rVB5	19163	47659	1.11%	0.111%
11	8.861	1171	1180	1189	rBV	580946	1167476	27.12%	2.715%
12	9.349	1249	1263	1268	rBV3	110443	320736	7.45%	0.746%
13	9.408	1268	1273	1281	rVV3	65809	133022	3.09%	0.309%
14	9.626	1299	1310	1323	rVB4	67011	163440	3.80%	0.380%
15	9.773	1327	1335	1343	rVB3	87279	182997	4.25%	0.426%
16	9.914	1349	1359	1363	rBV2	83509	178060	4.14%	0.414%
17	9.961	1363	1367	1373	rVB2	47264	89263	2.07%	0.208%
18	10.096	1383	1390	1394	rBV4	59341	123243	2.86%	0.287%
19	10.144	1394	1398	1406	rVB2	53506	94313	2.19%	0.219%
20	10.261	1414	1418	1423	rVB3	38426	76344	1.77%	0.178%
21	10.338	1423	1431	1446	rBV	1253443	2513352	58.38%	5.846%
22	10.455	1446	1451	1454	rBV3	41763	67735	1.57%	0.158%
23	10.502	1454	1459	1470	rVB4	127017	352784	8.19%	0.821%
24	10.632	1470	1481	1484	rBV4	61062	131776	3.06%	0.307%
25	10.679	1484	1489	1493	rVV2	136044	240196	5.58%	0.559%
26	10.779	1500	1506	1510	rBV4	99324	215543	5.01%	0.501%
27	10.814	1510	1512	1515	rVV3	40180	48748	1.13%	0.113%
28	10.861	1515	1520	1525	rVB	160521	261954	6.08%	0.609%
29	11.049	1541	1552	1560	rBV	272463	703528	16.34%	1.636%
30	11.120	1560	1564	1567	rVV2	120615	206838	4.80%	0.481%
31	11.191	1571	1576	1579	rVV	266733	472355	10.97%	1.099%
32	11.243	1579	1585	1591	rVV	796619	1537861	35.72%	3.577%
33	11.296	1591	1594	1597	rVV3	128500	201137	4.67%	0.468%
34	11.349	1597	1603	1608	rVV2	568418	1075411	24.98%	2.501%
35	11.391	1608	1610	1615	rVV	168538	224713	5.22%	0.523%
36	11.443	1615	1619	1627	rVB2	288705	524320	12.18%	1.220%

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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_D\Method\82D061121S.M
 Title : SW846 8260

37	11.514	1627	1631	1636	rBV	108921	175782	4.08%	0.409%
38	11.567	1636	1640	1644	rVV5	38865	65710	1.53%	0.153%
39	11.643	1645	1653	1663	rVB	958251	1761818	40.92%	4.098%
40	11.773	1670	1675	1678	rBV2	148171	231404	5.38%	0.538%
41	11.826	1678	1684	1689	rVV4	313377	703158	16.33%	1.636%
42	11.885	1689	1694	1697	rVV2	399161	714869	16.61%	1.663%
43	11.920	1698	1700	1706	rVB4	353675	552514	12.83%	1.285%
44	11.985	1706	1711	1715	rBV4	119564	233484	5.42%	0.543%
45	12.067	1718	1725	1731	rVB6	169510	389748	9.05%	0.907%
46	12.149	1731	1739	1745	rBV3	699482	1774847	41.23%	4.128%
47	12.232	1750	1753	1762	rVB6	220895	560451	13.02%	1.304%
48	12.373	1762	1777	1784	rBV4	1181581	4305109	100.00%	10.013%
49	12.455	1786	1791	1795	rBV6	210501	354497	8.23%	0.825%
50	12.514	1798	1801	1808	rVB5	336833	632821	14.70%	1.472%
51	12.573	1808	1811	1814	rVB2	115012	145557	3.38%	0.339%
52	12.626	1814	1820	1826	rVB2	1057123	1828147	42.46%	4.252%
53	12.714	1826	1835	1840	rBV6	304189	926614	21.52%	2.155%
54	12.790	1843	1848	1854	rVB2	909918	1675424	38.92%	3.897%
55	12.873	1854	1862	1867	rBV5	343586	1064250	24.72%	2.475%
56	12.955	1868	1876	1880	rBV8	550510	1507066	35.01%	3.505%
57	13.090	1894	1899	1903	rBV4	390254	596928	13.87%	1.388%
58	13.137	1903	1907	1910	rBV5	238527	415162	9.64%	0.966%
59	13.273	1926	1930	1932	rBV2	182157	252462	5.86%	0.587%
60	13.308	1932	1936	1939	rVV	266527	401596	9.33%	0.934%
61	13.355	1940	1944	1947	rVB3	249793	378229	8.79%	0.880%
62	13.432	1954	1957	1958	rBV3	107840	131541	3.06%	0.306%
63	13.490	1965	1967	1973	rVB6	159895	204298	4.75%	0.475%
64	13.567	1975	1980	1990	rVB	927108	1712631	39.78%	3.983%
65	13.708	1998	2004	2009	rBV3	398572	715137	16.61%	1.663%
66	13.749	2009	2011	2014	rVB4	62125	66166	1.54%	0.154%
67	13.890	2029	2035	2039	rBV2	611329	1094730	25.43%	2.546%
68	13.990	2049	2052	2056	rBV5	161907	258824	6.01%	0.602%
69	14.208	2085	2089	2092	rBV4	138585	203868	4.74%	0.474%
70	14.273	2097	2100	2107	rVB5	239838	439954	10.22%	1.023%
71	14.332	2108	2110	2115	rVB5	73019	88454	2.05%	0.206%
72	14.396	2116	2121	2124	rBV2	320493	509476	11.83%	1.185%
73	14.561	2141	2149	2154	rBV5	271405	637154	14.80%	1.482%
74	14.814	2189	2192	2197	rVB5	129845	206501	4.80%	0.480%
75	15.149	2245	2249	2251	rBV4	102659	154791	3.60%	0.360%
76	15.267	2266	2269	2278	rVB6	183234	358255	8.32%	0.833%

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ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
0228-COMPOSITE-E

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

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77	15.355	2280	2284	2291	rVB3	207920	350751	8.15%	0.816%
78	15.431	2293	2297	2300	rBV7	90373	157911	3.67%	0.367%
79	15.767	2351	2354	2361	rVB6	187465	360303	8.37%	0.838%

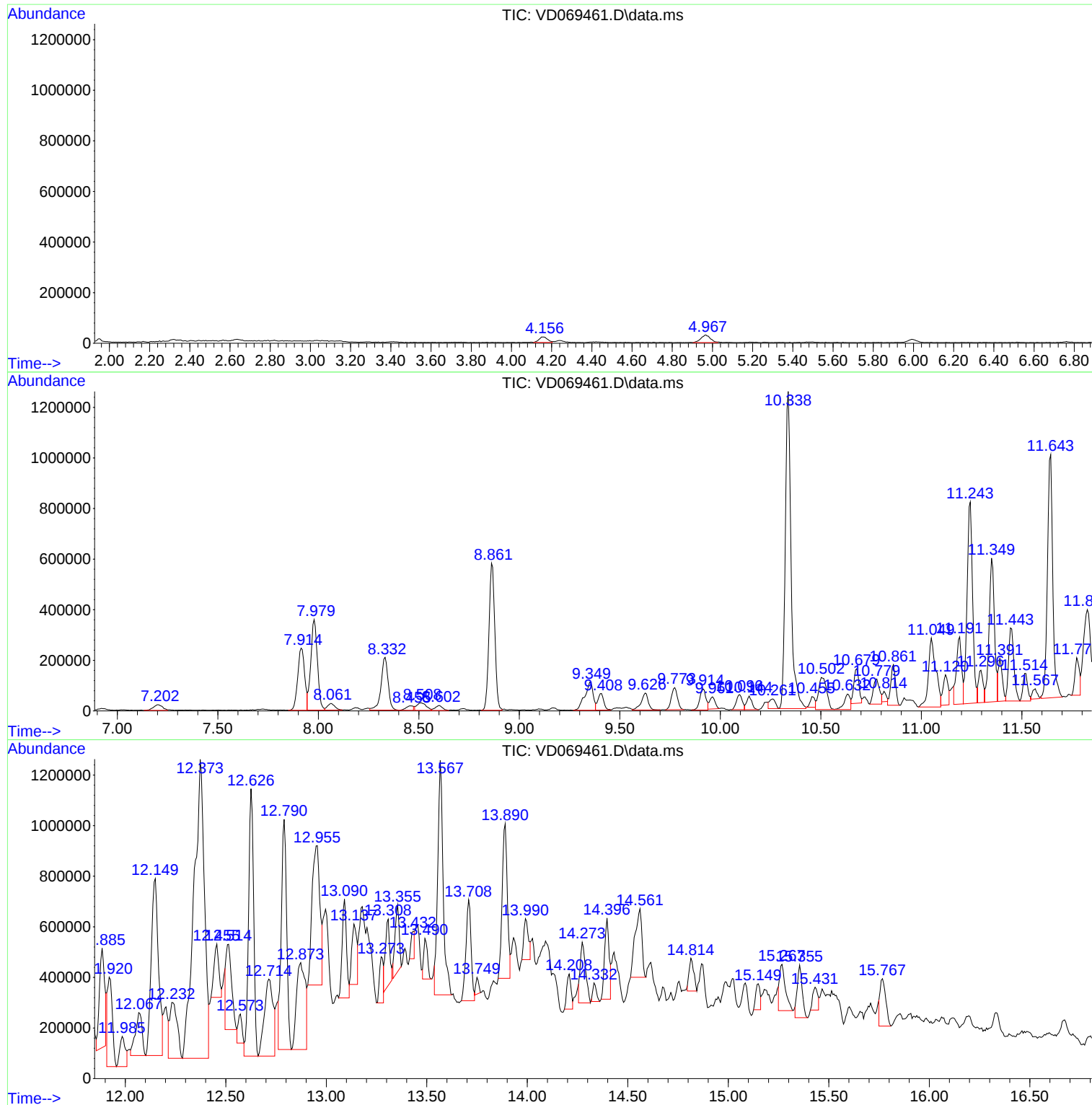
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Data File : VD069461.D
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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D061121S.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD061621\
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Operator : VA/SY
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Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

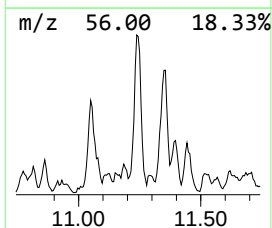
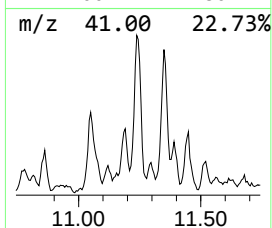
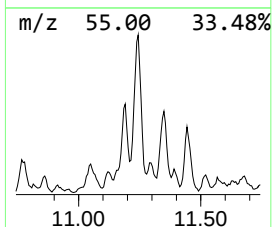
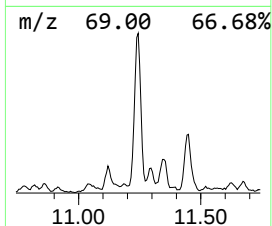
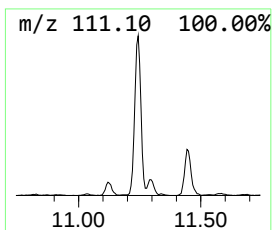
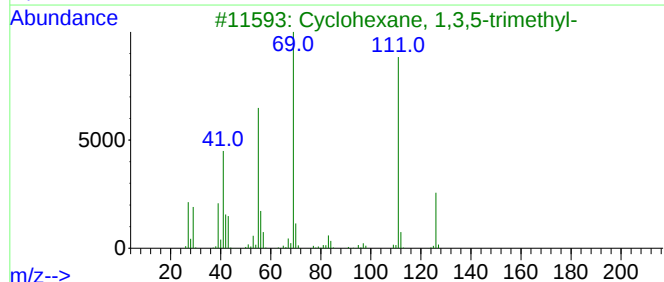
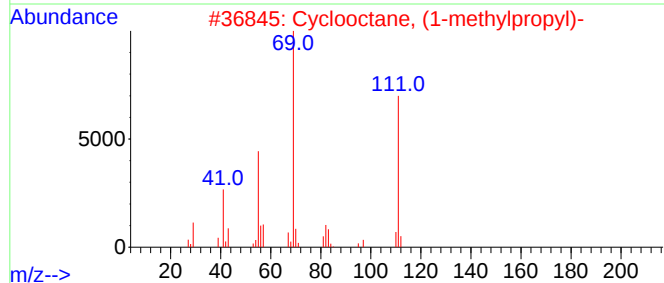
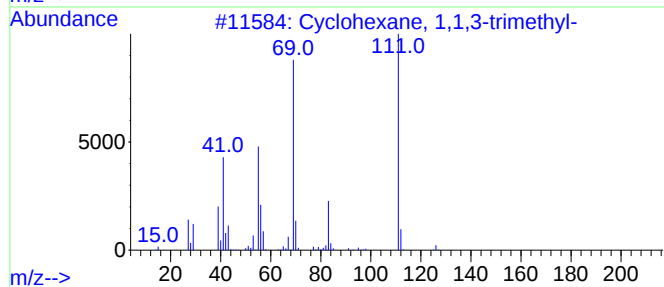
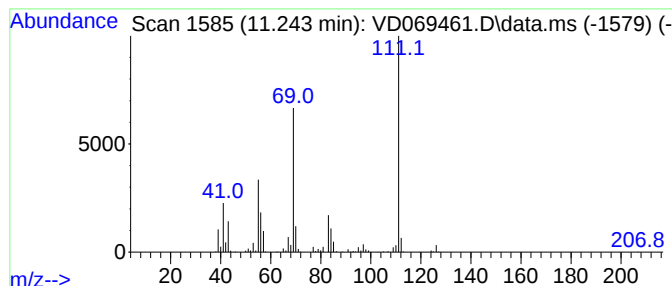
Instrument :
MSVOA_D
ClientSampleId :
0228-COMPOSITE-E

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D061121S.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.243	43.64 ug/l	1537860	Chlorobenzene-d5	11.643	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	72
2	Cyclooctane, (1-methylpropyl)-	168	C12H24	016538-89-9	64
3	Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	59
4	4-Amino-6-hydroxypyrimidine	111	C4H5N3O	001193-22-2	59
5	p-Fluoroaniline	111	C6H6FN	000371-40-4	58



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Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
0228-COMPOSITE-E

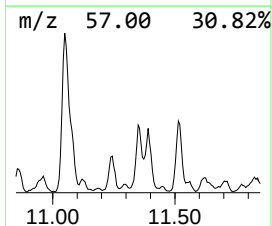
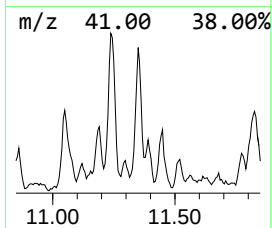
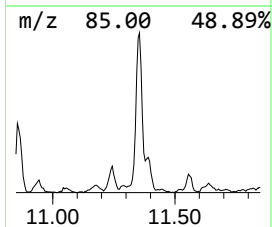
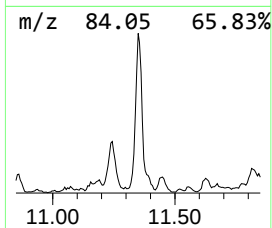
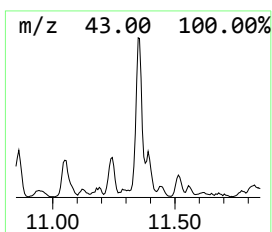
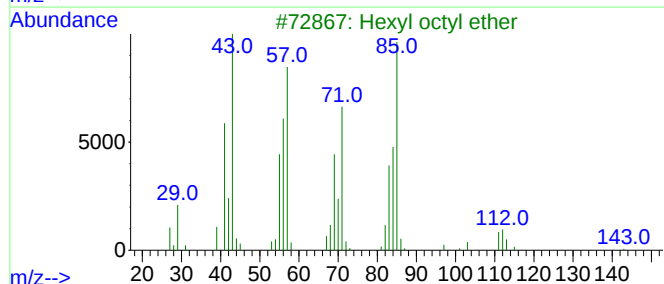
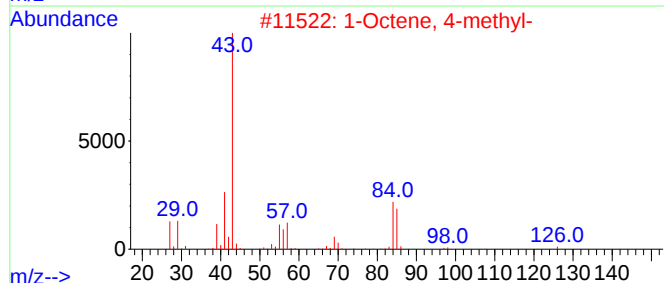
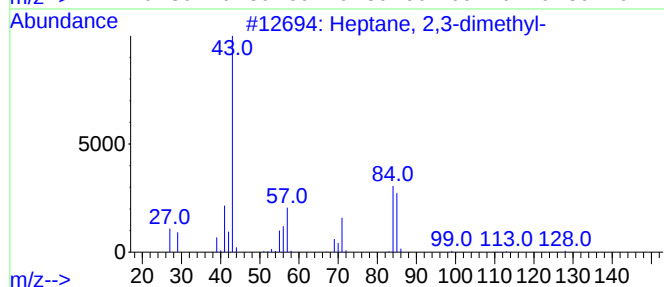
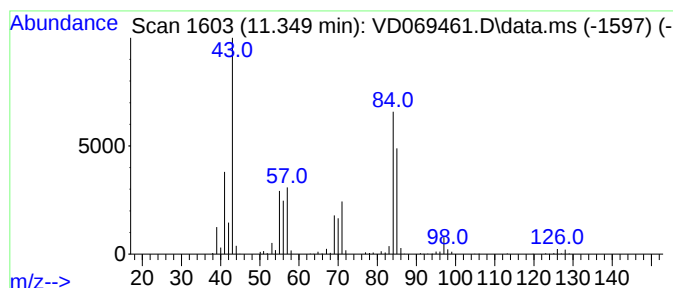
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D061121S.M
Quant Title : SW846 8260

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

Peak Number 2 Heptane, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.349	30.52 ug/l	1075410	Chlorobenzene-d5	11.643

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 2,3-dimethyl-	128	C9H20	003074-71-3	72
2		1-Octene, 4-methyl-	126	C9H18	013151-12-7	58
3		Hexyl octyl ether	214	C14H30O	017071-54-4	53
4		Hexane, 2,3,5-trimethyl-	128	C9H20	001069-53-0	53
5		Hexane, 3-ethyl-2-methyl-	128	C9H20	016789-46-1	50



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ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
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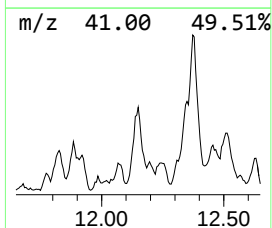
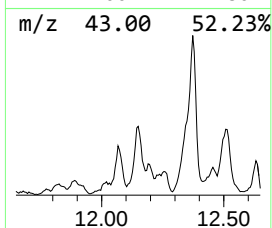
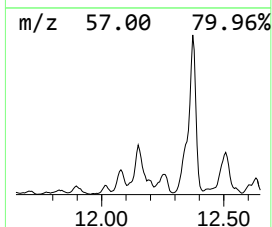
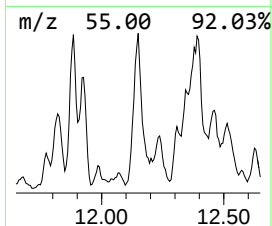
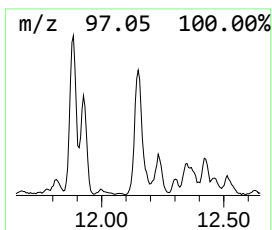
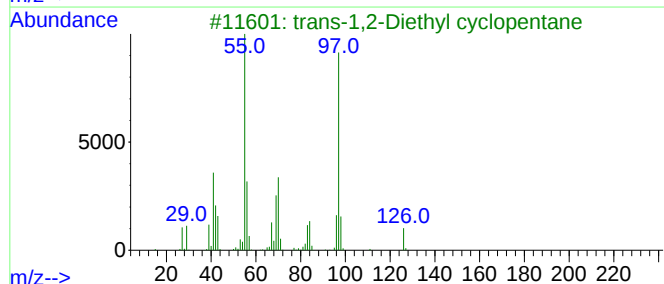
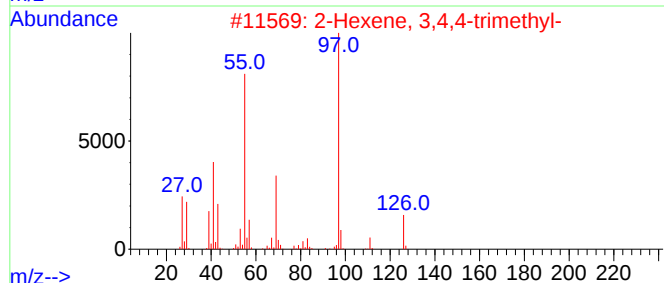
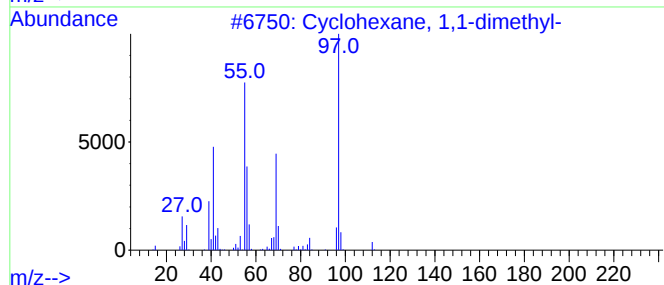
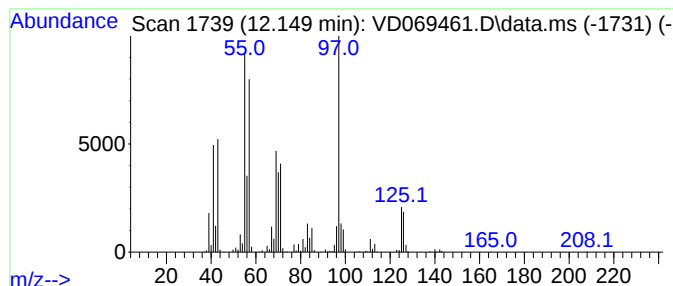
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D061121S.M
Quant Title : SW846 8260

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

Peak Number 3 unknown12.149 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.149	50.37 ug/l	1774850	Chlorobenzene-d5	11.643

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	43
2			2-Hexene, 3,4,4-trimethyl-	126	C9H18	053941-19-8	43
3			trans-1,2-Diethyl cyclopentane	126	C9H18	000932-40-1	43
4			Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	41
5			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	38



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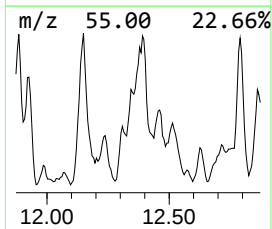
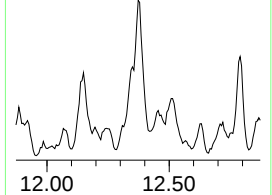
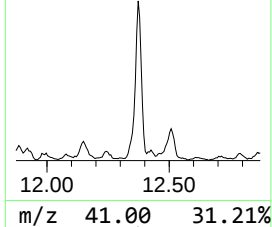
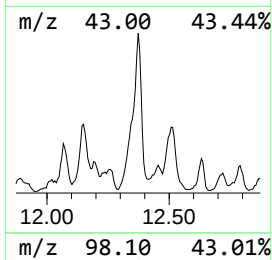
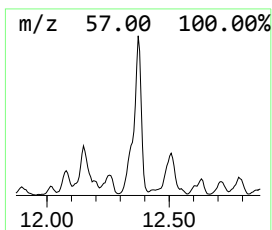
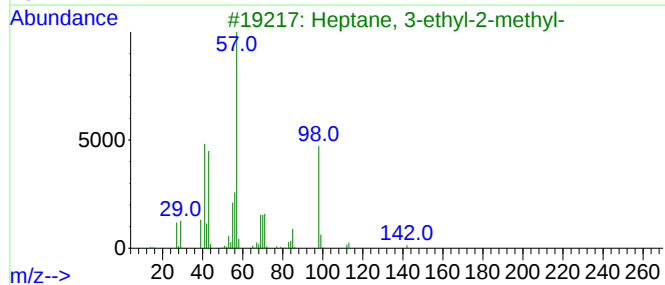
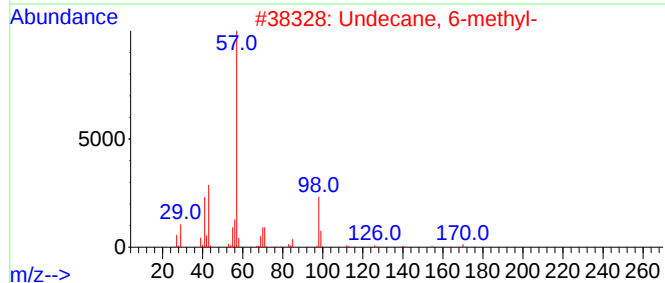
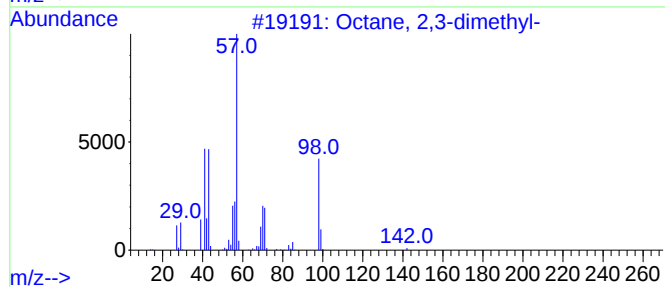
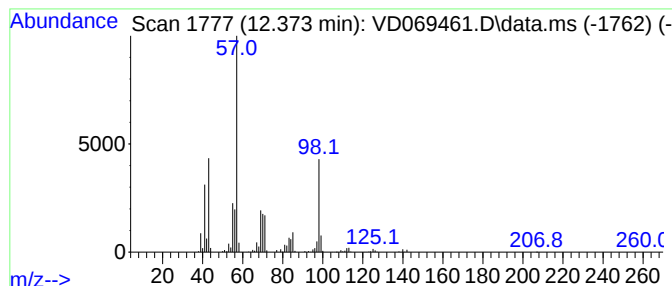
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Quant Title : SW846 8260

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

Peak Number 4 Octane, 2,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.373	122.18 ug/l	4305110	Chlorobenzene-d5	11.643

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	87
2			Undecane, 6-methyl-	170	C12H26	017302-33-9	59
3			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	58
4			Heptane, 3,5-dimethyl-	128	C9H20	000926-82-9	43
5			Heptyl isobutyl carbonate	216	C12H24O3	959068-08-7	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD061621\
Data File : VD069461.D
Acq On : 16 Jun 2021 12:32
Operator : VA/SY
Sample : M2709-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
0228-COMPOSITE-E

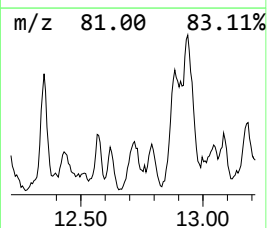
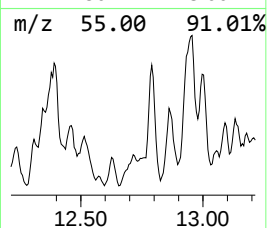
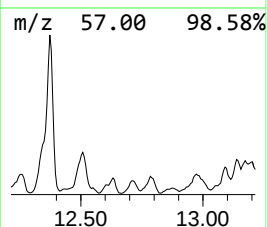
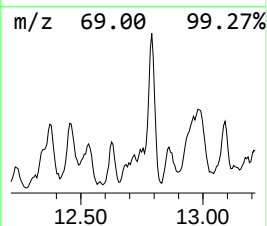
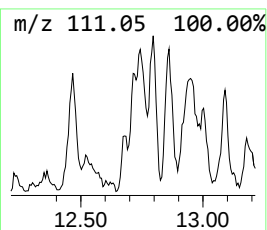
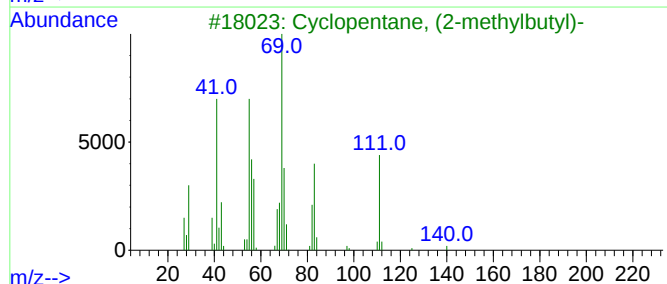
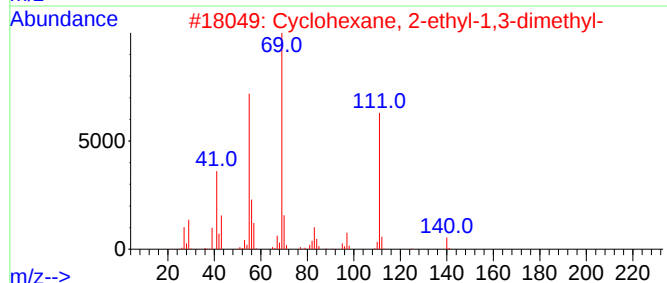
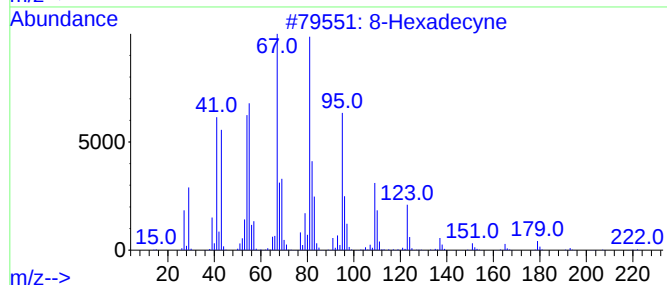
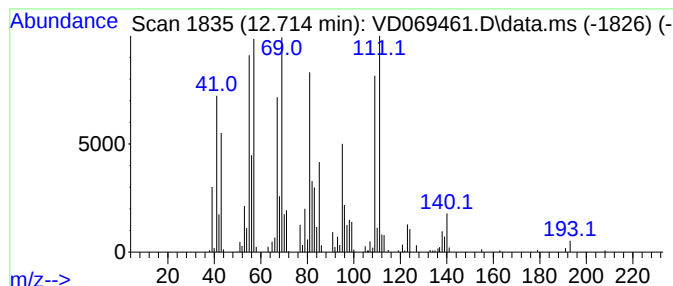
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Quant Title : SW846 8260

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

Peak Number 5 unknown12.714 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.714	27.05 ug/l	926614	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			8-Hexadecyne	222	C16H30	019781-86-3	25
2			Cyclohexane, 2-ethyl-1,3-dimethyl-	140	C10H20	007045-67-2	25
3			Cyclopentane, (2-methylbutyl)-	140	C10H20	053366-38-4	25
4			Bicyclo[4.1.0]heptane, 3-methyl-	110	C8H14	041977-47-3	25
5			Cyclohexane, 1-ethyl-2,3-dimethyl-	140	C10H20	007058-05-1	25



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD061621\
Data File : VD069461.D
Acq On : 16 Jun 2021 12:32
Operator : VA/SY
Sample : M2709-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
0228-COMPOSITE-E

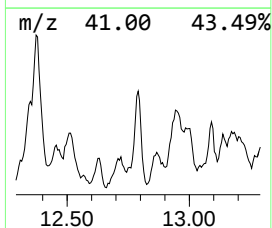
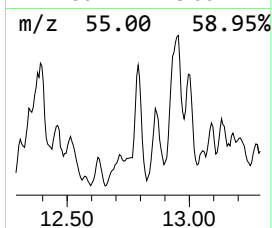
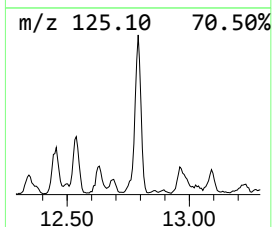
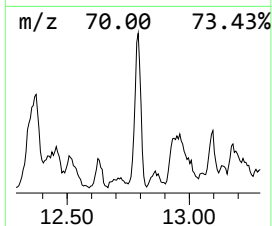
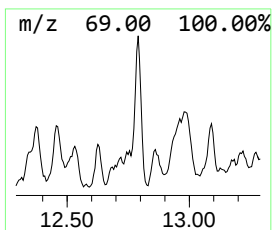
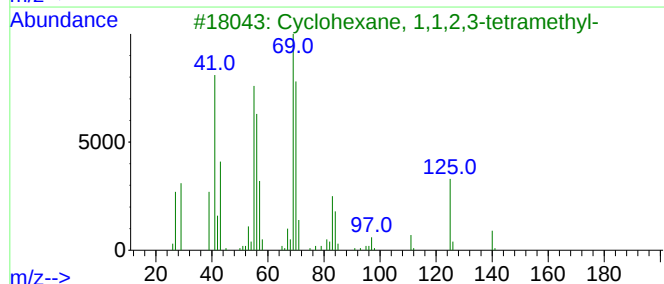
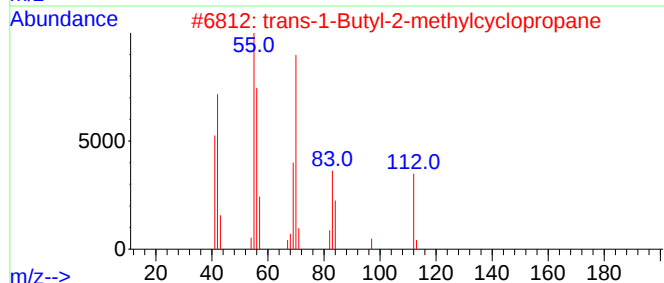
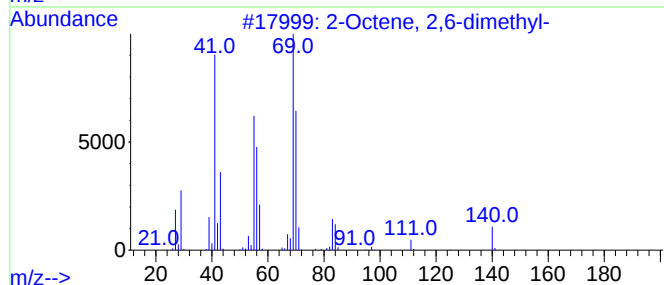
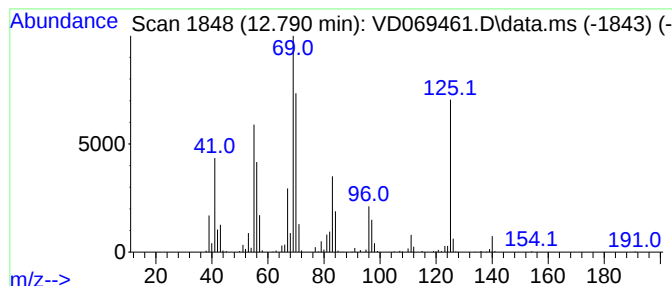
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Quant Title : SW846 8260

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

Peak Number 6 2-Octene, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.790	48.91 ug/l	1675420	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	70
2			trans-1-Butyl-2-methylcyclopropane	112	C8H16	038851-70-6	50
3			Cyclohexane, 1,1,2,3-tetramethyl-	140	C10H20	006783-92-2	46
4			Cyclooctane, 1-methyl-3-propyl-	168	C12H24	255885-37-1	38
5			cis-1-Butyl-2-methylcyclopropane	112	C8H16	038851-69-3	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD061621\
Data File : VD069461.D
Acq On : 16 Jun 2021 12:32
Operator : VA/SY
Sample : M2709-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
0228-COMPOSITE-E

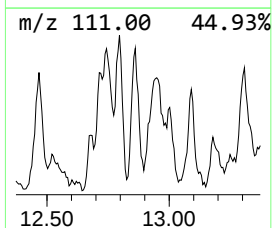
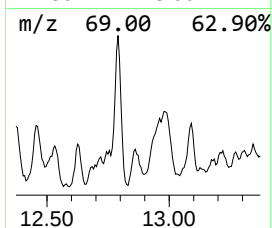
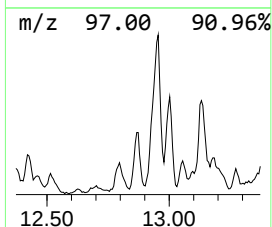
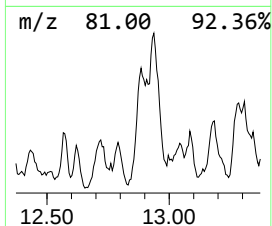
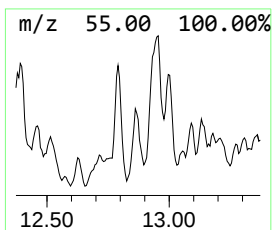
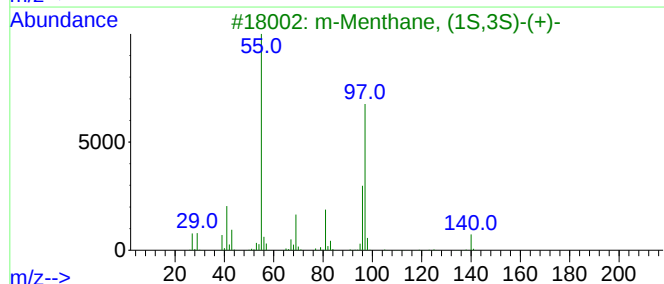
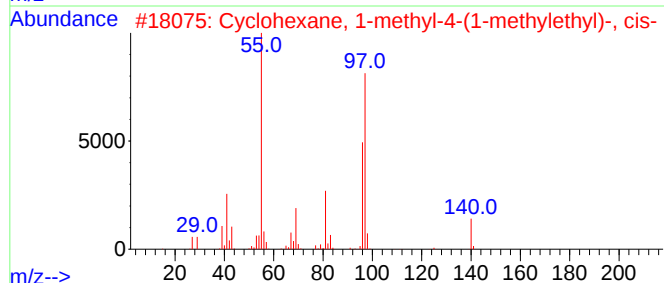
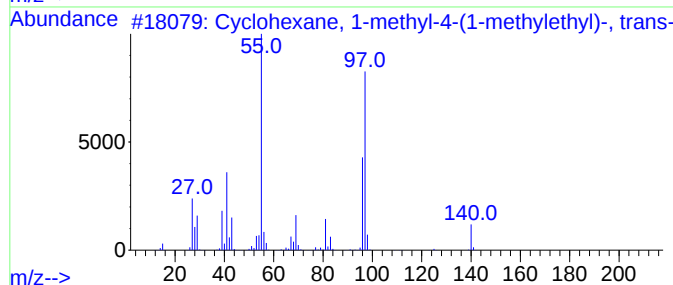
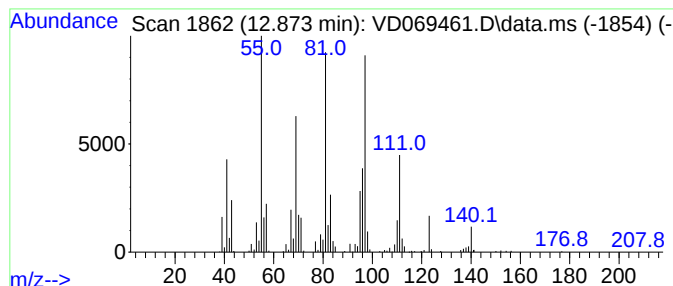
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Quant Title : SW846 8260

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

Peak Number 7 unknown12.873 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.873	31.07 ug/l	1064250	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	38
2			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	38
3			m-Menthane, (1S,3S)-(+)-	140	C10H20	013837-67-7	38
4			1-Methyl-4-(1-methylethyl)-cyclo...	140	C10H20	000099-82-1	35
5			Cyclohexane, 1-methyl-3-(1-methy...	140	C10H20	016580-24-8	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD061621\
Data File : VD069461.D
Acq On : 16 Jun 2021 12:32
Operator : VA/SY
Sample : M2709-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
0228-COMPOSITE-E

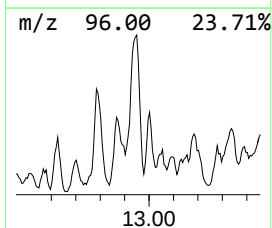
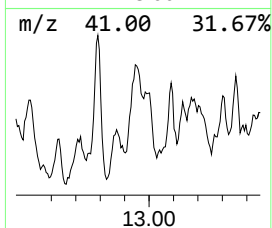
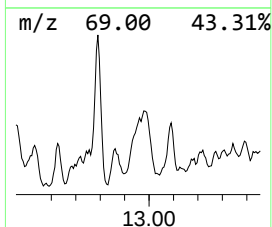
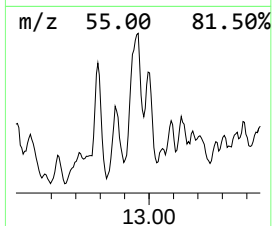
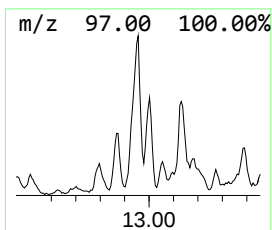
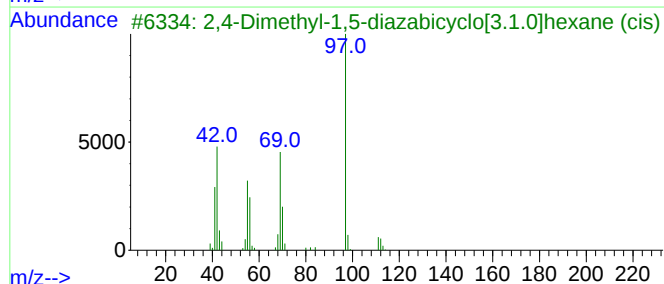
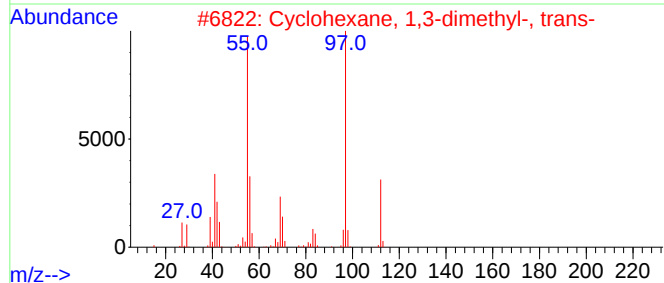
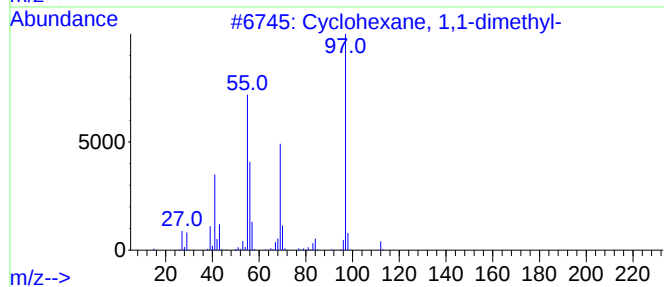
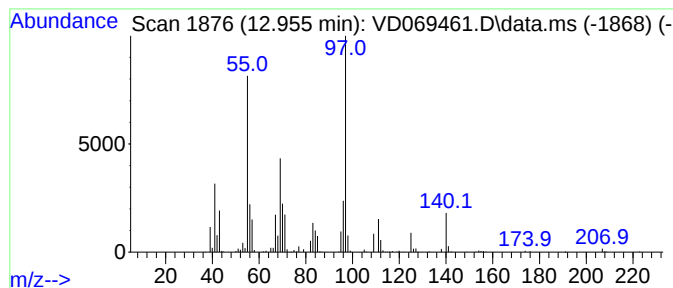
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Quant Title : SW846 8260

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

Peak Number 8 Cyclohexane, 1,1-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.955	44.00 ug/l	1507070	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	53
2			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	53
3			2,4-Dimethyl-1,5-diazabicyclo[3....	112	C6H12N2	100463-01-2	52
4			Oxalic acid, cyclohexylmethyl pr...	228	C12H20O4	1000309-68-1	47
5			4-Octene, 2,6-dimethyl-, [S-(E)]-	140	C10H20	062960-76-3	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD061621\
Data File : VD069461.D
Acq On : 16 Jun 2021 12:32
Operator : VA/SY
Sample : M2709-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
0228-COMPOSITE-E

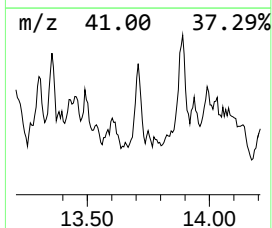
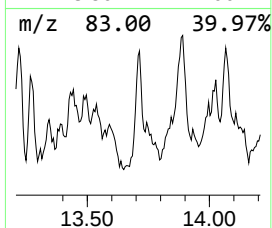
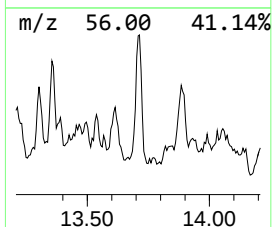
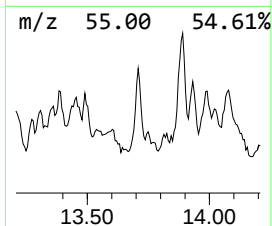
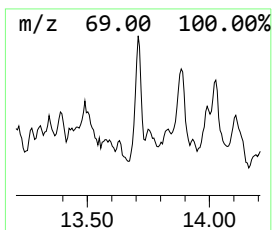
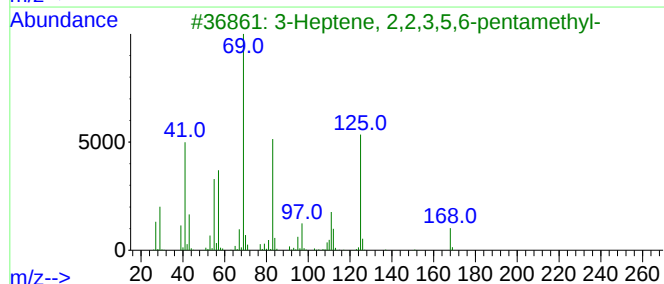
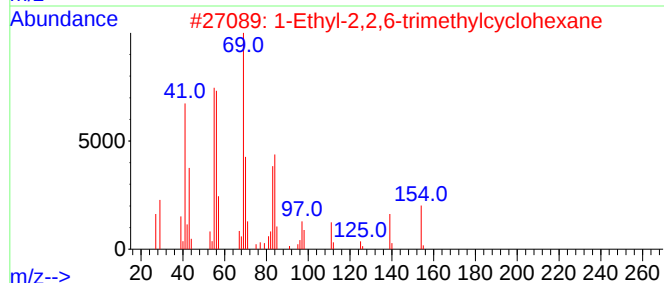
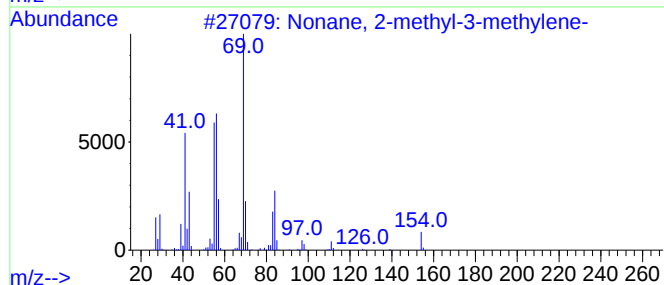
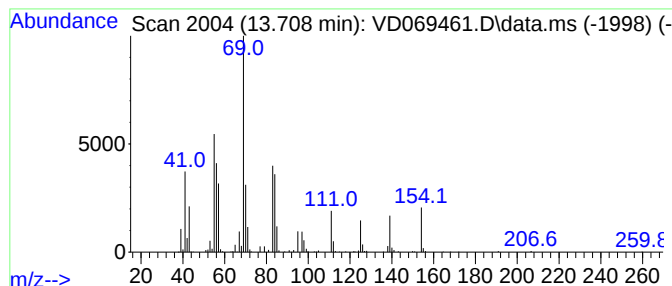
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Quant Title : SW846 8260

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

Peak Number 9 Nonane, 2-methyl-3-methylene- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.708	20.88 ug/l	715137	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 2-methyl-3-methylene-	154	C11H22	055499-08-6	68
2			1-Ethyl-2,2,6-trimethylcyclohexane	154	C11H22	071186-27-1	64
3			3-Heptene, 2,2,3,5,6-pentamethyl-	168	C12H24	116164-06-8	47
4			Cyclopentane, (2-methylbutyl)-	140	C10H20	053366-38-4	47
5			5-Undecene	154	C11H22	004941-53-1	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD061621\
Data File : VD069461.D
Acq On : 16 Jun 2021 12:32
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Sample : M2709-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
0228-COMPOSITE-E

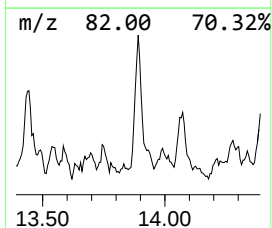
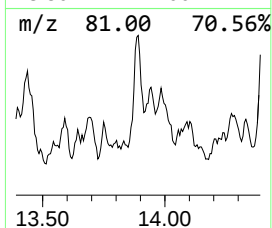
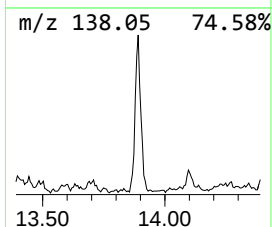
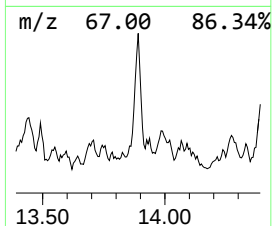
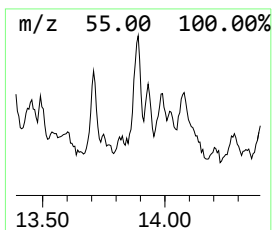
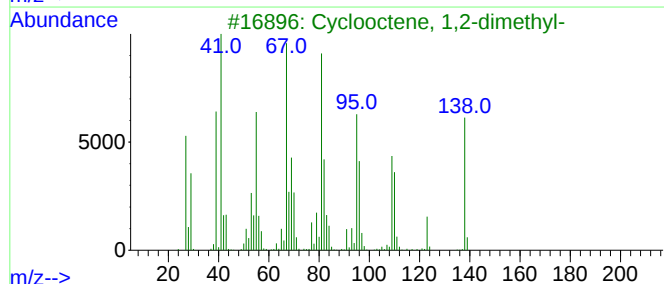
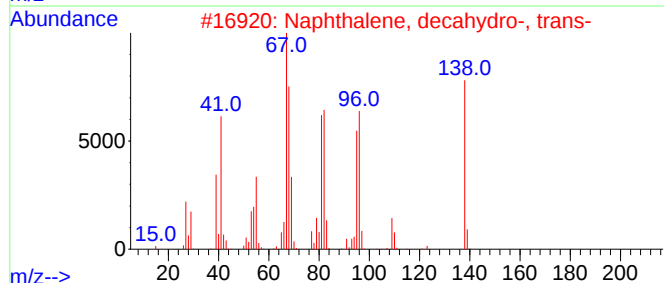
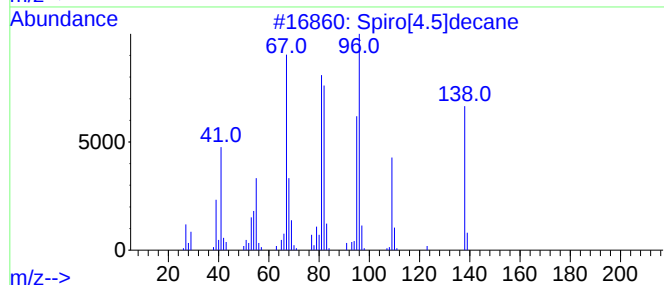
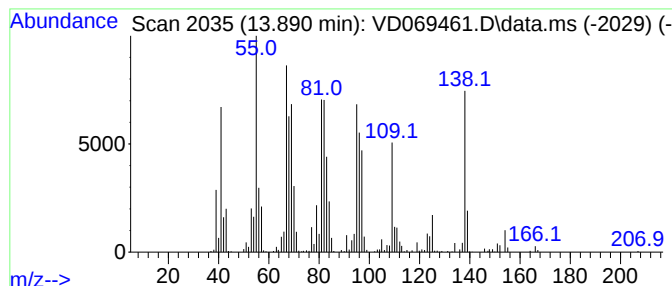
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Quant Title : SW846 8260

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

Peak Number 10 Spiro[4.5]decane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.890	31.96 ug/l	1094730	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Spiro[4.5]decane	138	C10H18	000176-63-6	70
2			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	70
3			Cyclooctene, 1,2-dimethyl-	138	C10H18	054299-96-6	68
4			1H-Inden-1-one, octahydro-, cis-	138	C9H14O	002826-65-5	62
5			9-Decen-1-ol, trifluoroacetate	252	C12H19F3O2	1000352-65-9	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD061621\
Data File : VD069461.D
Acq On : 16 Jun 2021 12:32
Operator : VA/SY
Sample : M2709-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
0228-COMPOSITE-E

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D061121S.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclohexane, 1,...	11.243	43.6	ug/l	1537860	3	11.643	1761820	50.0
Heptane, 2,3-di...	11.349	30.5	ug/l	1075410	3	11.643	1761820	50.0
unknown12.149	12.149	50.4	ug/l	1774850	3	11.643	1761820	50.0
Octane, 2,3-dim...	12.373	122.2	ug/l	4305110	3	11.643	1761820	50.0
unknown12.714	12.714	27.1	ug/l	926614	4	13.567	1712630	50.0
2-Octene, 2,6-d...	12.790	48.9	ug/l	1675420	4	13.567	1712630	50.0
unknown12.873	12.873	31.1	ug/l	1064250	4	13.567	1712630	50.0
Cyclohexane, 1,...	12.955	44.0	ug/l	1507070	4	13.567	1712630	50.0
Nonane, 2-methy...	13.708	20.9	ug/l	715137	4	13.567	1712630	50.0
Spiro[4.5]decane	13.890	32.0	ug/l	1094730	4	13.567	1712630	50.0