

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121121\
 Data File : VD071277.D
 Acq On : 11 Dec 2021 18:41
 Operator : VA/SY
 Sample : M5051-03
 Misc : 5.04G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 P-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_D\Method\82D112921S.M
 Title : SW846 8260

Signal : TIC: VD071277.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.962	506	517	533	rBV3	57583	200481	7.69%	0.562%
2	7.185	887	895	904	rBV2	12737	37085	1.42%	0.104%
3	7.914	1009	1019	1024	rBV	220007	543994	20.86%	1.526%
4	7.979	1024	1030	1040	rVB	442257	988464	37.91%	2.773%
5	8.326	1074	1089	1100	rVB2	190522	592668	22.73%	1.663%
6	8.861	1172	1180	1189	rBV	659287	1304168	50.02%	3.659%
7	10.338	1421	1431	1439	rBV	1004464	1856386	71.20%	5.208%
8	10.397	1439	1441	1452	rVB2	26292	42882	1.64%	0.120%
9	10.720	1490	1496	1501	rVB2	37554	59744	2.29%	0.168%
10	11.079	1546	1557	1560	rBV7	11408	29762	1.14%	0.084%
11	11.238	1579	1584	1590	rVB3	31961	59275	2.27%	0.166%
12	11.349	1598	1603	1607	rBV4	17787	30754	1.18%	0.086%
13	11.449	1607	1620	1627	rVB8	68211	180063	6.91%	0.505%
14	11.526	1627	1633	1638	rBV3	29192	50791	1.95%	0.143%
15	11.638	1644	1652	1661	rBV	1061544	1818995	69.77%	5.104%
16	11.738	1663	1669	1678	rVV	189178	344742	13.22%	0.967%
17	11.843	1678	1687	1698	rVV	575193	1283754	49.24%	3.602%
18	11.920	1698	1700	1706	rVB3	47989	73925	2.84%	0.207%
19	11.979	1706	1710	1713	rBV4	19309	33307	1.28%	0.093%
20	12.073	1722	1726	1730	rVB4	27780	48239	1.85%	0.135%
21	12.173	1730	1743	1750	rBV	486268	1096273	42.05%	3.076%
22	12.261	1750	1758	1762	rVB2	132461	238171	9.14%	0.668%
23	12.373	1762	1777	1784	rBV4	232016	885583	33.97%	2.485%
24	12.473	1788	1794	1798	rBV5	109932	240350	9.22%	0.674%
25	12.549	1804	1807	1810	rBV2	49237	52387	2.01%	0.147%
26	12.620	1814	1819	1829	rVB	944697	1600088	61.37%	4.489%
27	12.708	1829	1834	1838	rBV6	52161	105742	4.06%	0.297%
28	12.785	1842	1847	1855	rVB3	258658	584656	22.42%	1.640%
29	12.873	1855	1862	1865	rBV	679213	1161364	44.54%	3.258%
30	12.908	1865	1868	1870	rVV	359257	554401	21.26%	1.555%
31	12.949	1870	1875	1881	rVV3	1358414	2607216	100.00%	7.315%
32	12.996	1881	1883	1888	rVB3	117267	164840	6.32%	0.462%
33	13.108	1894	1902	1909	rBV	432153	999147	38.32%	2.803%
34	13.173	1909	1913	1921	rVB3	394500	647341	24.83%	1.816%
35	13.255	1921	1927	1933	rBV	1312809	2098669	80.49%	5.888%
36	13.302	1933	1935	1939	rVB3	72106	79360	3.04%	0.223%

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Integration Parameters: RTEINT.P

Integrator: RTE

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Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
 Title : SW846 8260

37	13.349	1940	1943	1946	rBV3	55573	81350	3.12%	0.228%
38	13.385	1946	1949	1953	rVB4	98212	135505	5.20%	0.380%
39	13.455	1954	1961	1965	rBV4	367351	776290	29.77%	2.178%
40	13.496	1965	1968	1972	rVV3	187314	270515	10.38%	0.759%
41	13.561	1974	1979	1982	rVV	1140117	1813632	69.56%	5.088%
42	13.590	1982	1984	1989	rVB	798207	1117184	42.85%	3.134%
43	13.632	1989	1991	1998	rVB4	112593	145568	5.58%	0.408%
44	13.726	2000	2007	2009	rBV4	157902	331074	12.70%	0.929%
45	13.761	2009	2013	2017	rVV2	728175	1163420	44.62%	3.264%
46	13.796	2017	2019	2028	rVB2	379909	563916	21.63%	1.582%
47	13.885	2028	2034	2039	rBV2	706583	1174967	45.07%	3.297%
48	13.955	2039	2046	2050	rBV3	104003	204833	7.86%	0.575%
49	14.020	2053	2057	2059	rBV2	139252	193817	7.43%	0.544%
50	14.102	2066	2071	2076	rVB	326153	494067	18.95%	1.386%
51	14.214	2084	2090	2094	rVB2	254741	421931	16.18%	1.184%
52	14.273	2095	2100	2106	rVV8	71763	178291	6.84%	0.500%
53	14.349	2107	2113	2116	rVV3	123572	253516	9.72%	0.711%
54	14.396	2116	2121	2124	rVV3	228827	442615	16.98%	1.242%
55	14.432	2124	2127	2129	rVV	172158	256782	9.85%	0.720%
56	14.467	2129	2133	2137	rVV	265231	410847	15.76%	1.153%
57	14.508	2137	2140	2144	rVV2	103874	145347	5.57%	0.408%
58	14.555	2144	2148	2155	rVB5	149562	273855	10.50%	0.768%
59	14.649	2161	2164	2168	rVB	67400	87255	3.35%	0.245%
60	14.696	2168	2172	2175	rBV2	44891	74382	2.85%	0.209%
61	14.737	2175	2179	2184	rVB3	109025	180979	6.94%	0.508%
62	14.802	2184	2190	2197	rBV2	374584	770938	29.57%	2.163%
63	14.967	2214	2218	2224	rVB2	60878	93298	3.58%	0.262%
64	15.032	2225	2229	2233	rBV7	18304	30704	1.18%	0.086%
65	15.079	2233	2237	2240	rBV5	35963	53208	2.04%	0.149%
66	15.184	2250	2255	2261	rVB5	45599	104719	4.02%	0.294%
67	15.384	2277	2289	2295	rBV	299999	571072	21.90%	1.602%
68	15.479	2301	2305	2311	rVB7	14259	29121	1.12%	0.082%
69	16.473	2467	2474	2483	rVB2	30038	74401	2.85%	0.209%
70	16.679	2501	2509	2510	rBV2	17320	27542	1.06%	0.077%

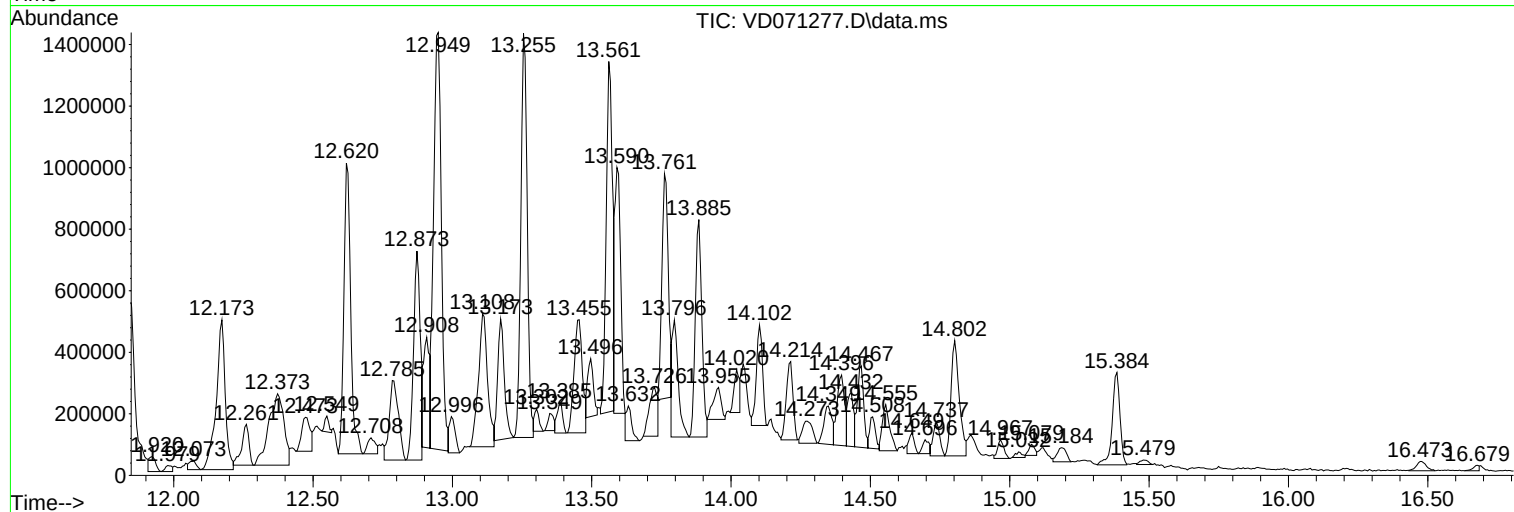
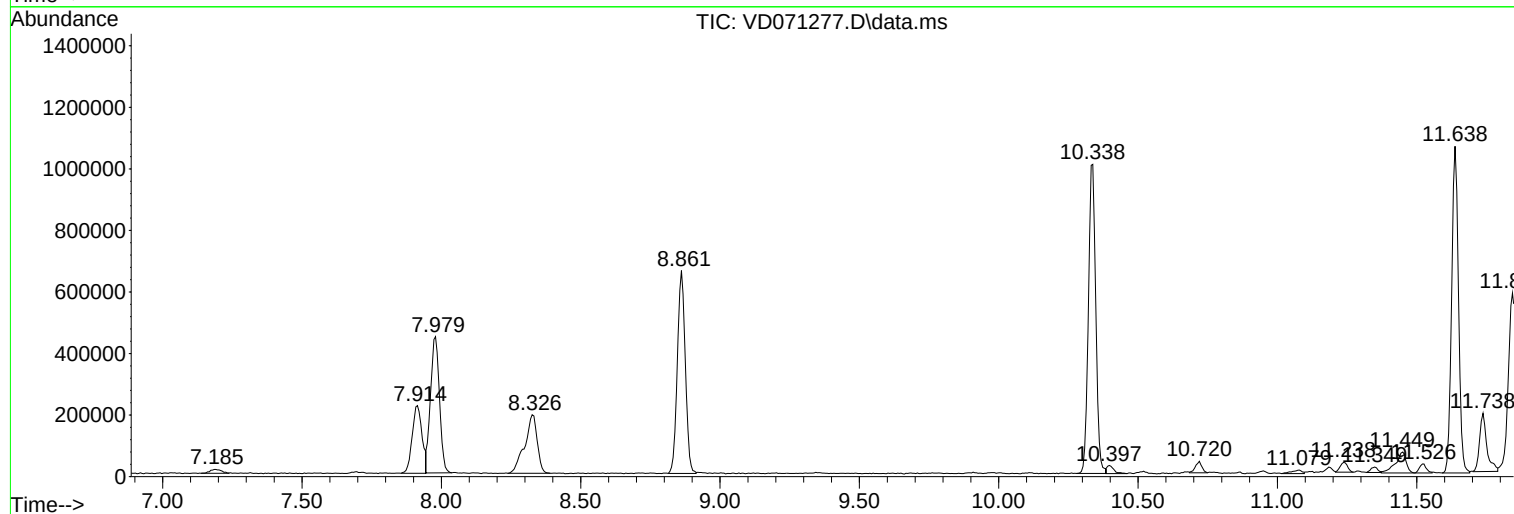
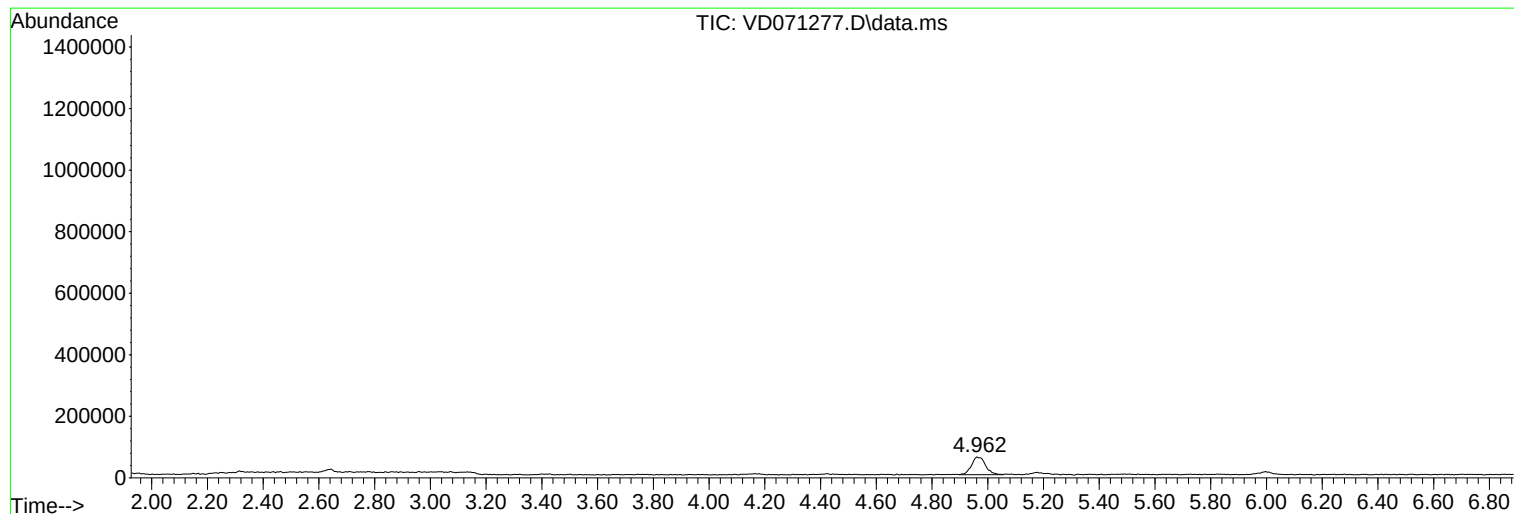
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Instrument :
MSVOA_D
ClientSampleId :
P-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121121\
Data File : VD071277.D
Acq On : 11 Dec 2021 18:41
Operator : VA/SY
Sample : M5051-03
Misc : 5.04G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
P-3

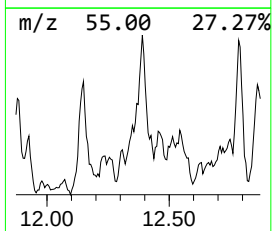
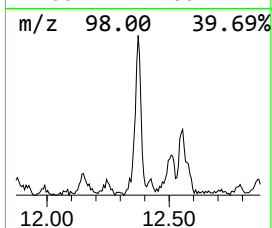
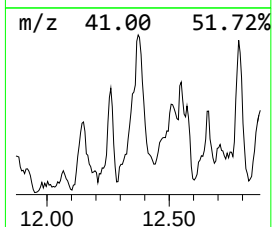
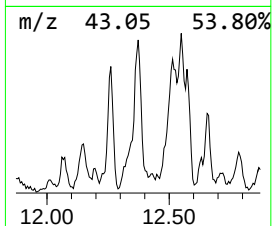
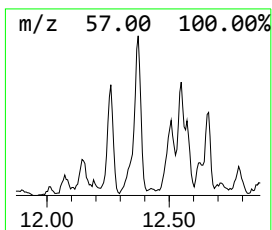
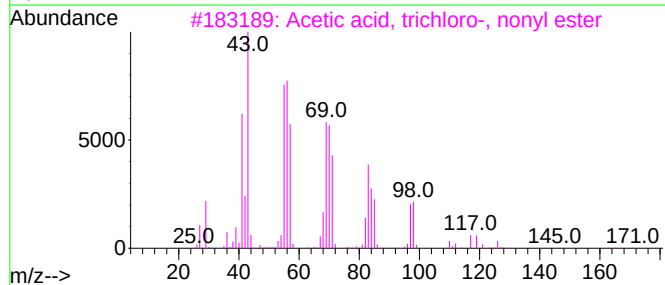
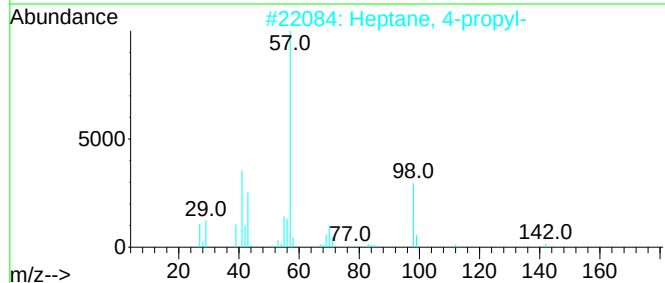
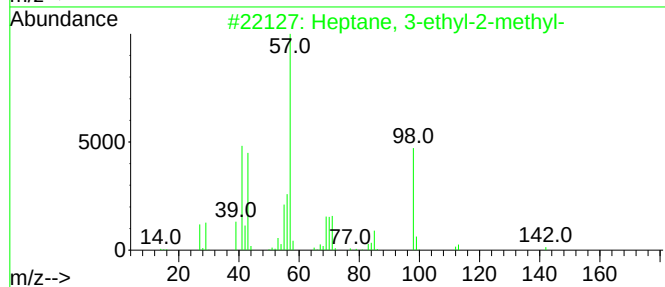
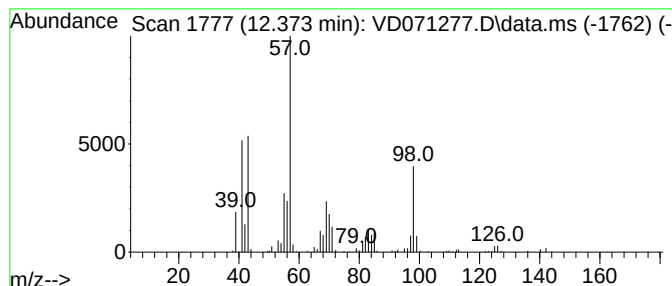
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Heptane, 3-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.373	24.34 ug/l	885583	Chlorobenzene-d5	11.638

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	83
2			Heptane, 4-propyl-	142	C10H22	003178-29-8	59
3			Acetic acid, trichloro-, nonyl e...	288	C11H19Cl3O2	065611-32-7	49
4			2- Chloropropionic acid, decyl e...	248	C13H25ClO2	086711-78-6	43
5			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	43



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Instrument :
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P-3

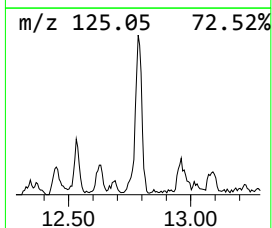
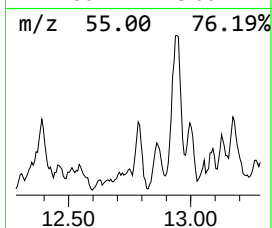
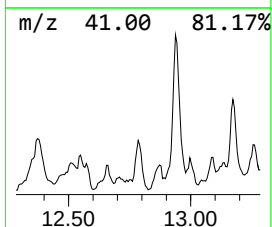
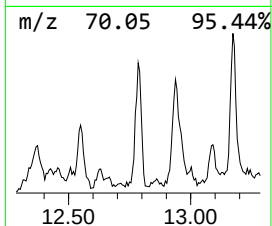
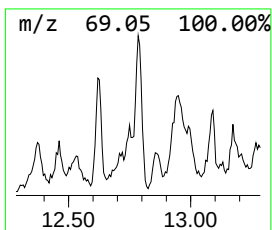
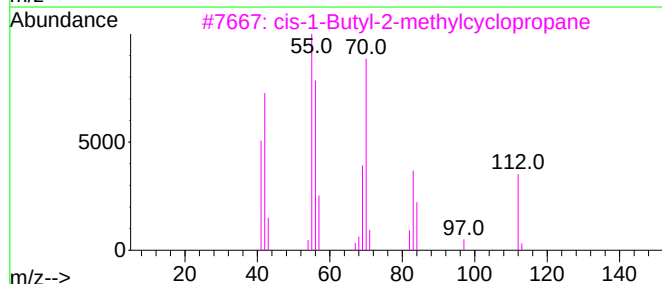
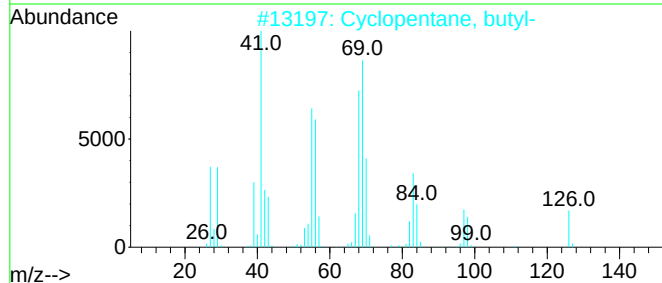
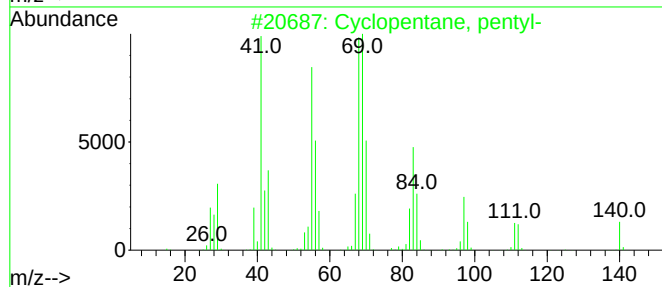
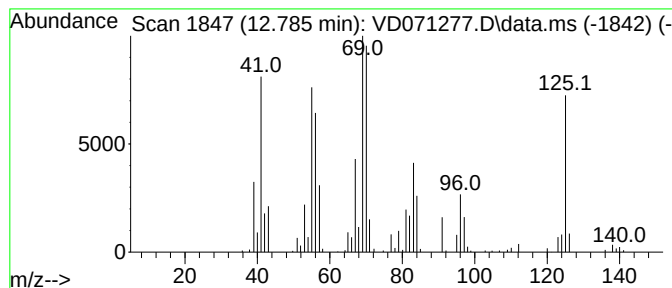
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclopentane, pentyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.785	16.12 ug/l	584656	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, pentyl-	140	C10H20	003741-00-2	58
2			Cyclopentane, butyl-	126	C9H18	002040-95-1	53
3			cis-1-Butyl-2-methylcyclopropane	112	C8H16	038851-69-3	50
4			2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	49
5			trans-1-Butyl-2-methylcyclopropane	112	C8H16	038851-70-6	45



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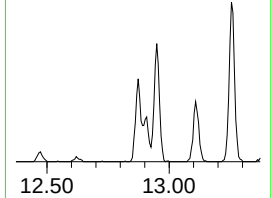
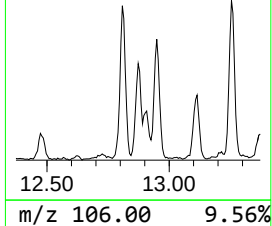
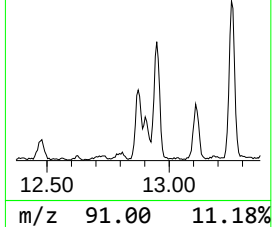
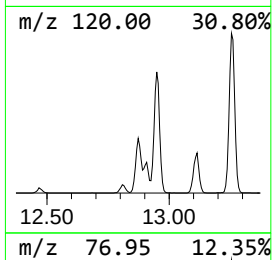
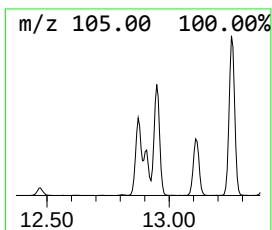
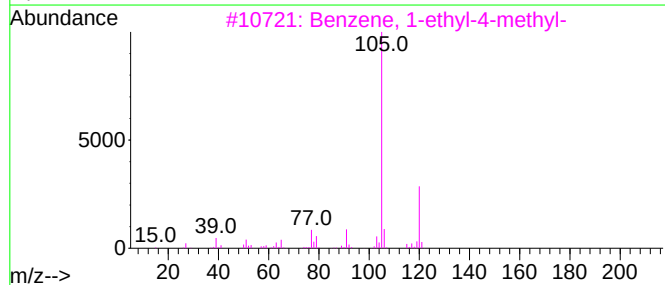
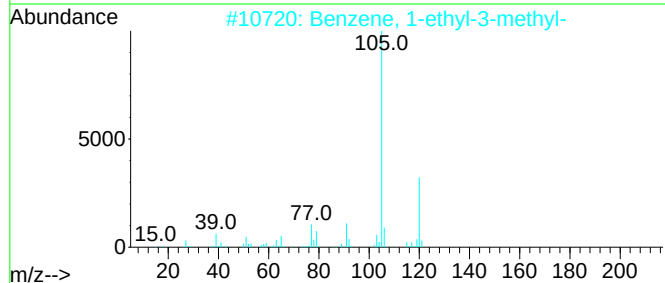
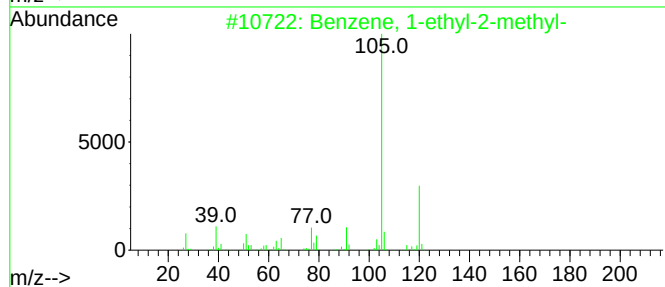
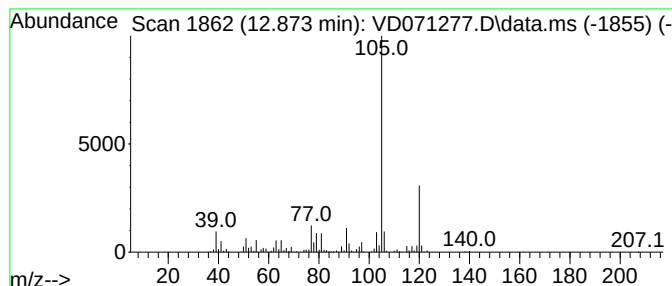
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.873	32.02 ug/l	1161360	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
5			Mesitylene	120	C9H12	000108-67-8	90



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Misc : 5.04G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
P-3

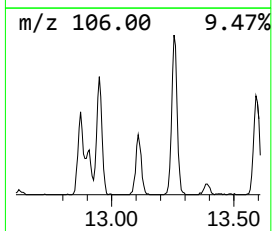
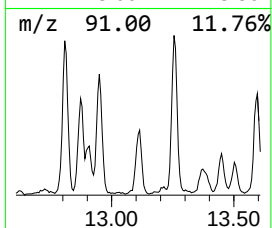
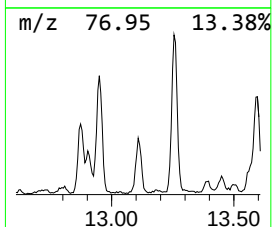
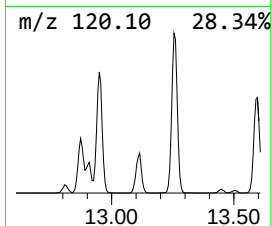
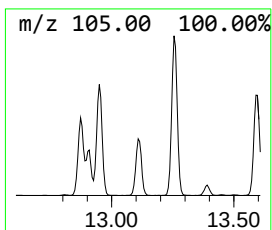
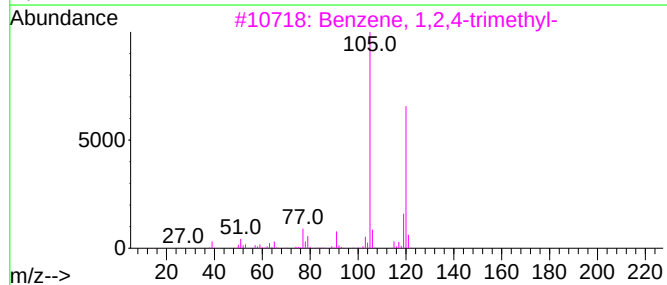
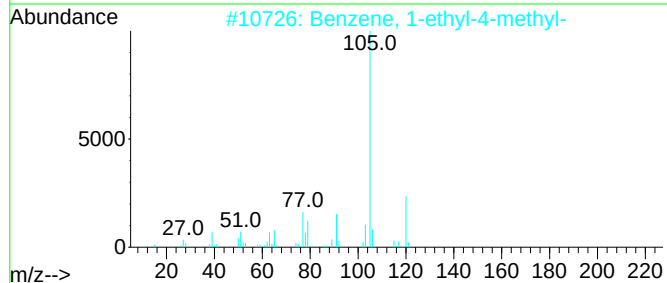
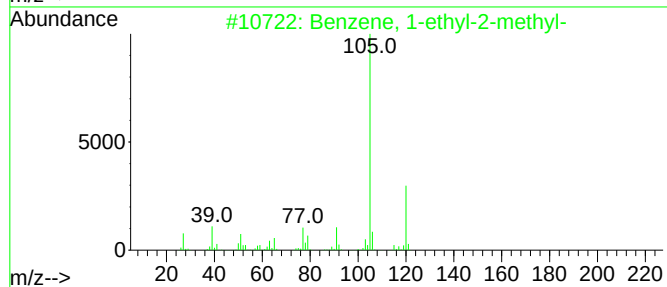
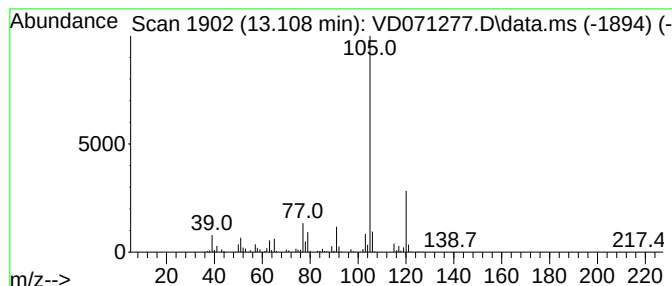
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1-ethyl-4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.108	27.55 ug/l	999147	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121121\
Data File : VD071277.D
Acq On : 11 Dec 2021 18:41
Operator : VA/SY
Sample : M5051-03
Misc : 5.04G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
P-3

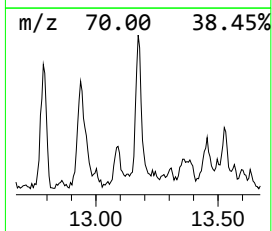
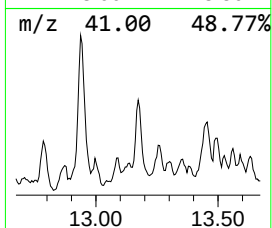
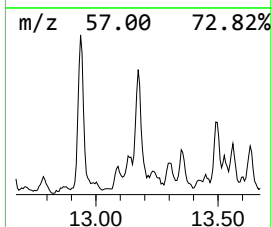
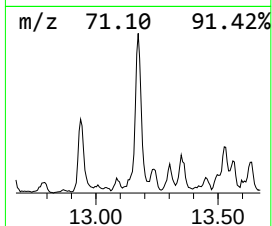
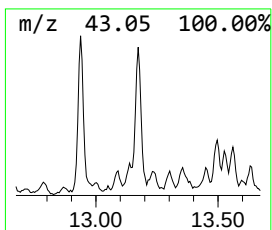
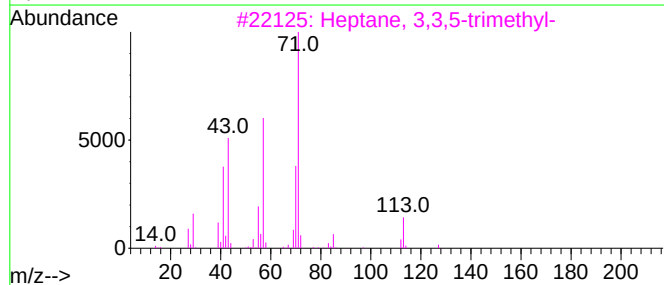
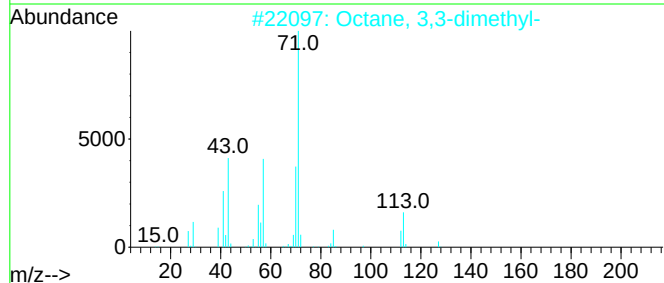
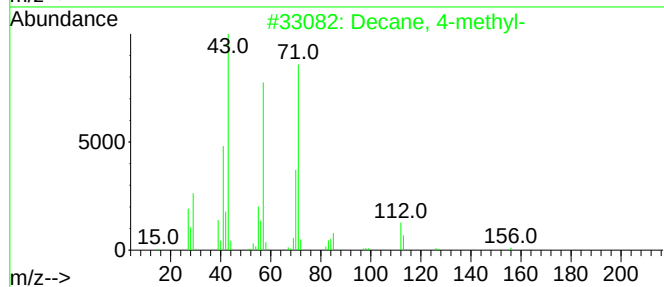
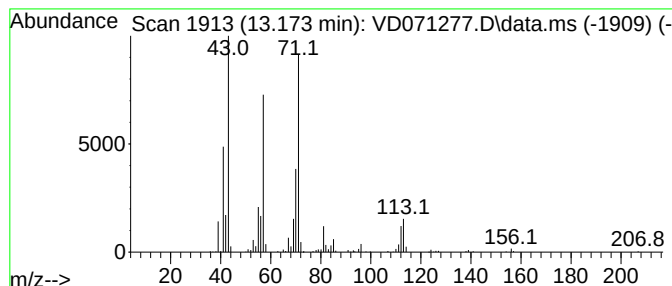
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Decane, 4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.173	17.85 ug/l	647341	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 4-methyl-	156	C11H24	002847-72-5	74
2			Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	72
3			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	64
4			Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	64
5			3,4-Dimethylpent-2-en-1-ol	114	C7H14O	1000195-06-9	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121121\
Data File : VD071277.D
Acq On : 11 Dec 2021 18:41
Operator : VA/SY
Sample : M5051-03
Misc : 5.04G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
P-3

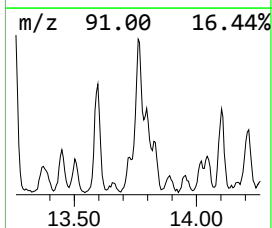
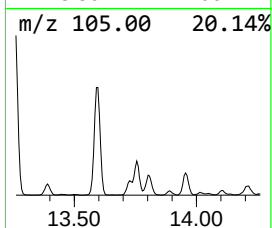
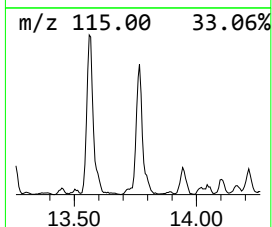
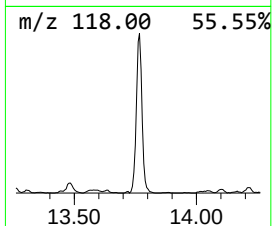
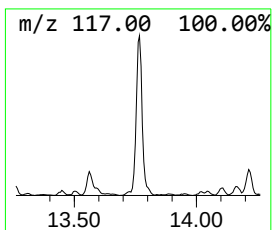
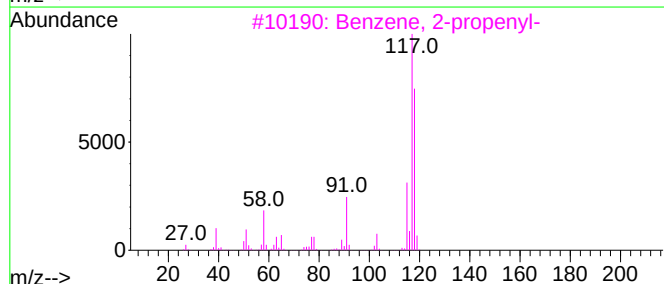
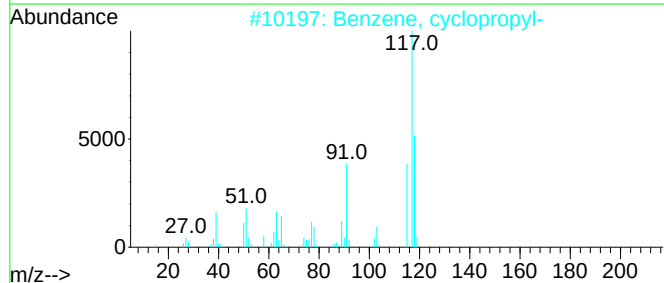
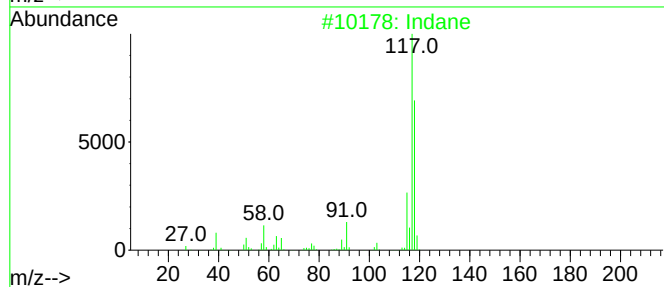
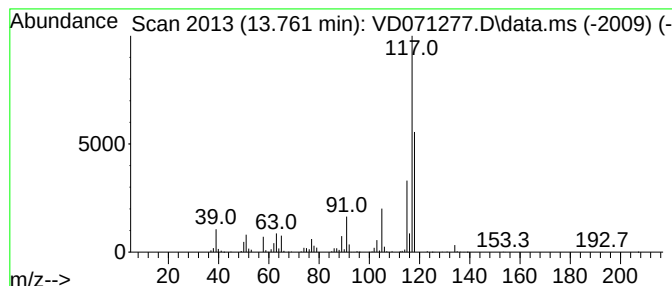
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.761	32.07 ug/l	1163420	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	81
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	68
4			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	62
5			Benzene, 1-propenyl-	118	C9H10	000637-50-3	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121121\
Data File : VD071277.D
Acq On : 11 Dec 2021 18:41
Operator : VA/SY
Sample : M5051-03
Misc : 5.04G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
P-3

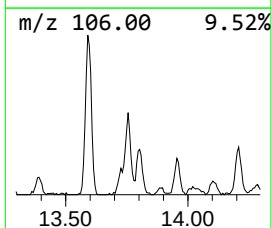
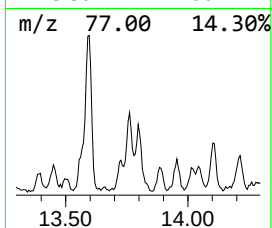
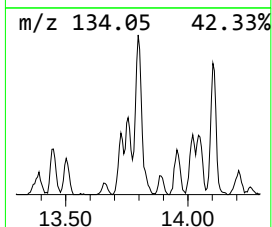
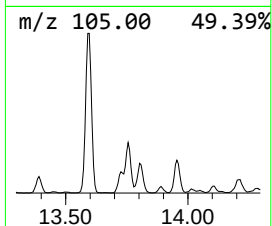
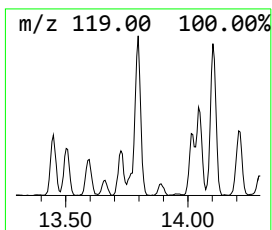
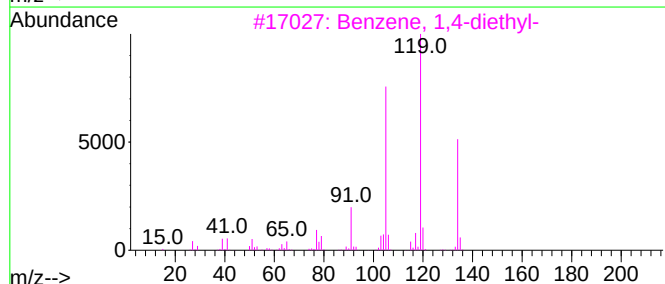
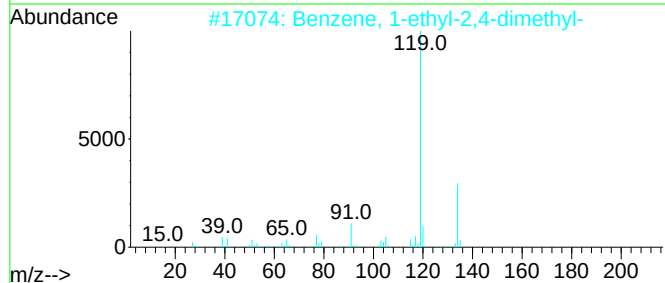
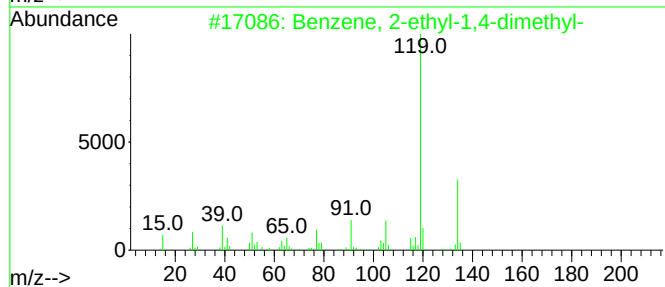
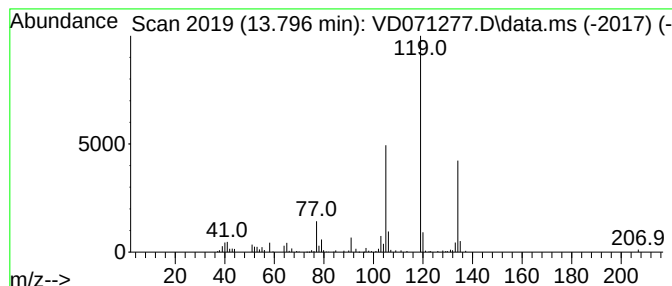
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.796	15.55 ug/l	563916	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	90
2			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	86
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	86
4			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	86
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121121\
Data File : VD071277.D
Acq On : 11 Dec 2021 18:41
Operator : VA/SY
Sample : M5051-03
Misc : 5.04G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
P-3

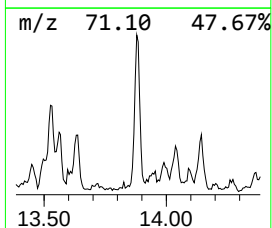
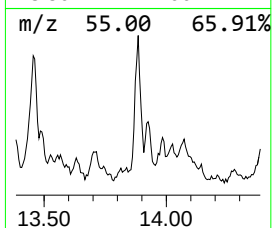
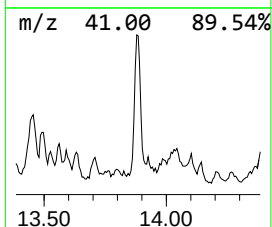
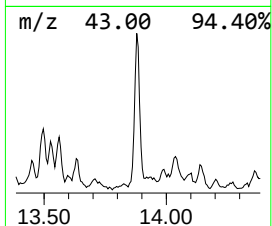
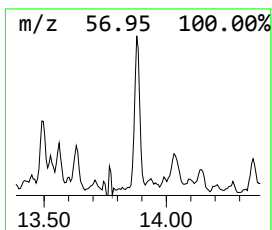
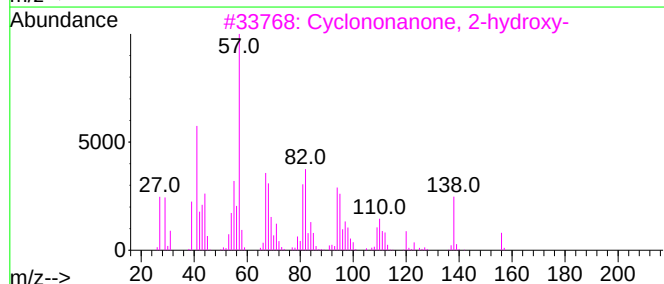
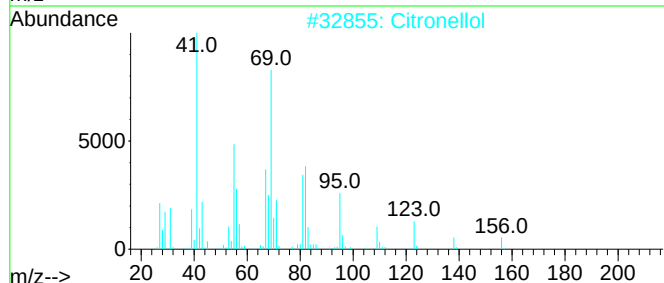
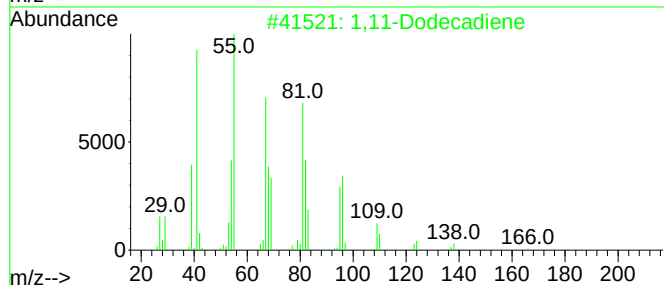
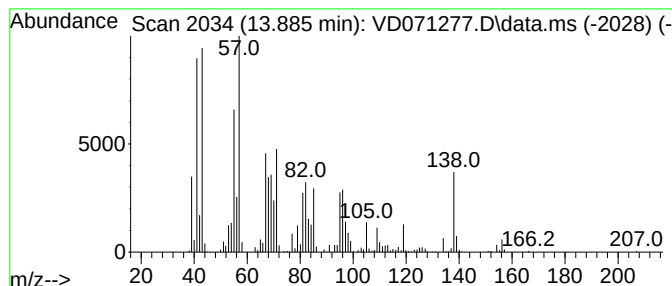
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1,11-Dodecadiene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.885	32.39 ug/l	1174970	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,11-Dodecadiene	166	C12H22	005876-87-9	50
2			Citronellol	156	C10H20O	000106-22-9	43
3			Cyclononane, 2-hydroxy-	156	C9H16O2	000496-83-3	38
4			9-Borabicyclo[3.3.1]nonane, 9-hy...	138	C8H15BO	063366-65-4	30
5			2-Propanone, 1-cyclohexylidene-	138	C9H14O	000874-68-0	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121121\
Data File : VD071277.D
Acq On : 11 Dec 2021 18:41
Operator : VA/SY
Sample : M5051-03
Misc : 5.04G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
P-3

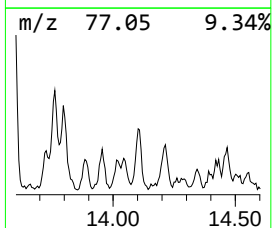
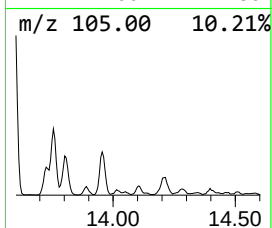
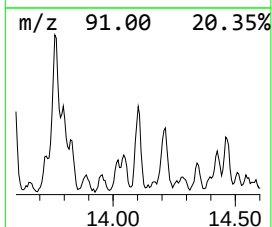
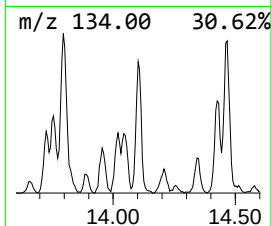
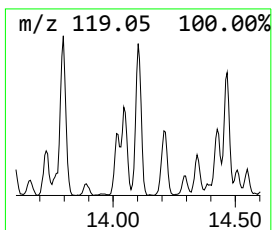
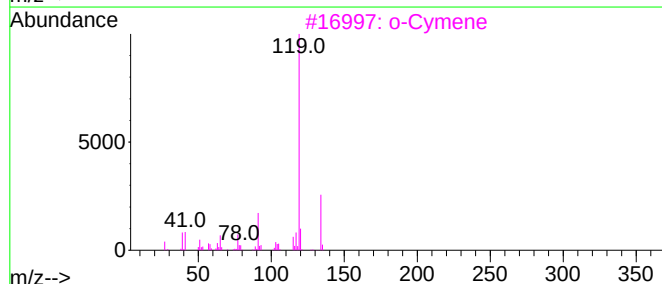
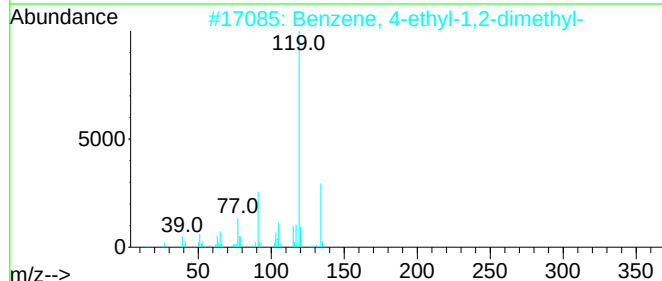
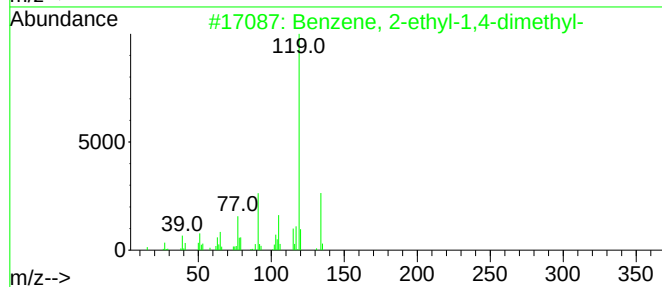
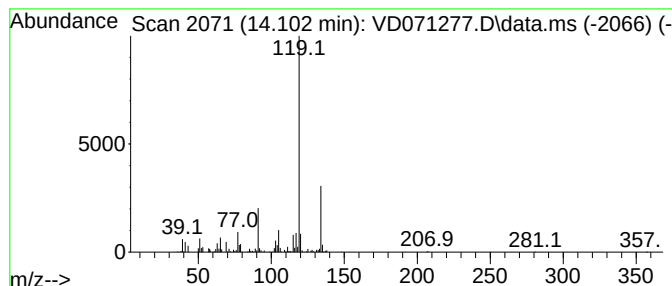
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.102	13.62 ug/l	494067	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5			p-Cymene	134	C10H14	000099-87-6	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121121\
Data File : VD071277.D
Acq On : 11 Dec 2021 18:41
Operator : VA/SY
Sample : M5051-03
Misc : 5.04G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

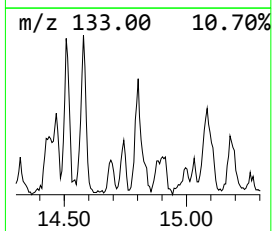
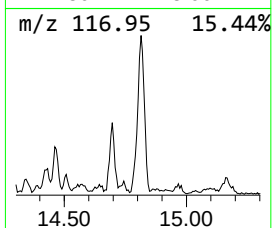
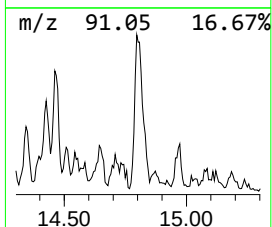
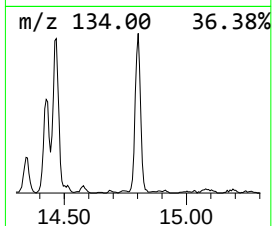
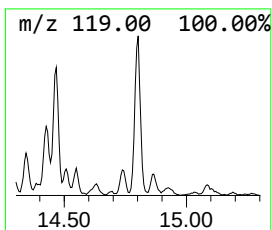
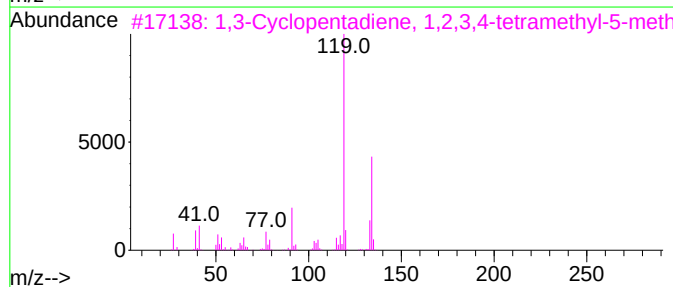
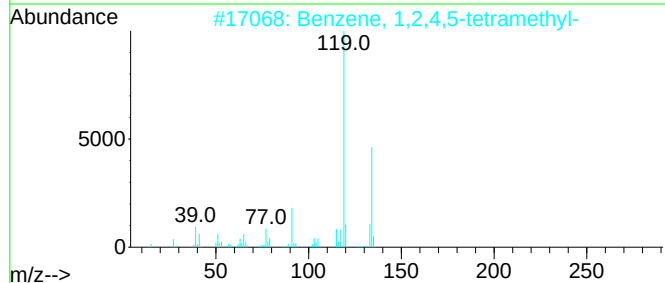
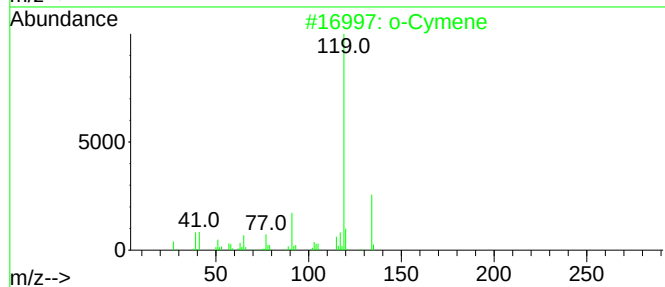
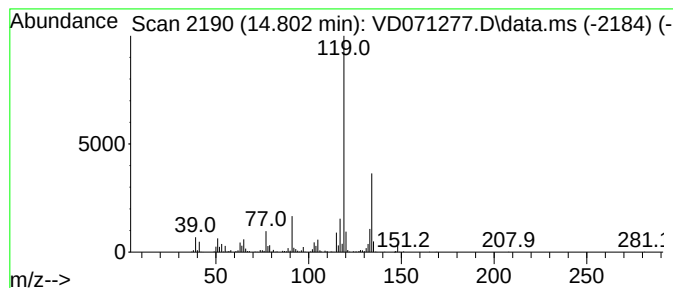
Instrument :
MSVOA_D
ClientSampleId :
P-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.802	21.25 ug/l	770938	1,4-Dichlorobenzene-d4			13.561
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	o-Cymene		134	C10H14	000527-84-4	95
2	Benzene, 1,2,4,5-tetramethyl-		134	C10H14	000095-93-2	94
3	1,3-Cyclopentadiene, 1,2,3,4-tet...		134	C10H14	076089-59-3	93
4	Benzene, 1,2,3,5-tetramethyl-		134	C10H14	000527-53-7	93
5	Benzene, 1-ethyl-2,4-dimethyl-		134	C10H14	000874-41-9	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121121\
Data File : VD071277.D
Acq On : 11 Dec 2021 18:41
Operator : VA/SY
Sample : M5051-03
Misc : 5.04G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
P-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Cyclopentane, p...	12.785	16.1	ug/l	584656	4	13.561	1813630	50.0
Benzene, 1-ethy...	12.873	32.0	ug/l	1161360	4	13.561	1813630	50.0
Benzene, 1-ethy...	13.108	27.6	ug/l	999147	4	13.561	1813630	50.0
Decane, 4-methyl-	13.173	17.9	ug/l	647341	4	13.561	1813630	50.0
Indane	13.761	32.1	ug/l	1163420	4	13.561	1813630	50.0
Benzene, 2-ethy...	13.796	15.6	ug/l	563916	4	13.561	1813630	50.0
1,11-Dodecadiene	13.885	32.4	ug/l	1174970	4	13.561	1813630	50.0
Benzene, 4-ethy...	14.102	13.6	ug/l	494067	4	13.561	1813630	50.0
o-Cymene	14.802	21.3	ug/l	770938	4	13.561	1813630	50.0