

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022522\
 Data File : VD072106.D
 Acq On : 25 Feb 2022 17:58
 Operator : VA/SY
 Sample : N1655-06
 Misc : 4.43G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 PREM-REAR

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\82D021822S.M
 Title : SW846 8260

Signal : TIC: VD072106.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.168	34	42	47	rBV	153648	207138	4.70%	0.346%
2	2.344	62	72	80	rBV	288435	457622	10.39%	0.764%
3	3.038	180	190	204	rBV	329169	717699	16.29%	1.199%
4	3.409	243	253	267	rVB2	190080	486202	11.03%	0.812%
5	3.879	324	333	344	rBV2	60953	167460	3.80%	0.280%
6	4.150	366	379	390	rBV3	29780	107483	2.44%	0.179%
7	4.950	504	515	517	rBV3	73300	177233	4.02%	0.296%
8	5.015	519	526	544	rVB3	166668	656847	14.91%	1.097%
9	5.491	591	607	621	rBV3	124855	435385	9.88%	0.727%
10	5.997	682	693	706	rBV2	134565	408748	9.28%	0.683%
11	6.344	744	752	767	rVB5	35101	113998	2.59%	0.190%
12	6.779	814	826	837	rBV5	28542	93020	2.11%	0.155%
13	6.926	840	851	860	rBV2	52048	151218	3.43%	0.253%
14	7.026	860	868	879	rVV2	119533	340098	7.72%	0.568%
15	7.732	981	988	997	rVB3	42060	109126	2.48%	0.182%
16	7.914	1009	1019	1023	rBV	468456	1130608	25.66%	1.888%
17	7.973	1023	1029	1038	rVV2	1020893	2627940	59.64%	4.389%
18	8.061	1038	1044	1055	rVB2	126273	332097	7.54%	0.555%
19	8.185	1057	1065	1073	rVV	277178	638443	14.49%	1.066%
20	8.344	1081	1092	1103	rVV3	799517	2260837	51.31%	3.776%
21	8.461	1103	1112	1113	rVV5	114218	231167	5.25%	0.386%
22	8.508	1114	1120	1130	rVV	303155	914958	20.77%	1.528%
23	8.597	1130	1135	1145	rVV4	91435	230129	5.22%	0.384%
24	8.720	1145	1156	1166	rVB	409943	856659	19.44%	1.431%
25	8.861	1171	1180	1192	rVV	1425737	2983576	67.71%	4.983%
26	9.091	1213	1219	1223	rBV3	29678	65370	1.48%	0.109%
27	9.138	1223	1227	1237	rVB4	31464	82022	1.86%	0.137%
28	9.355	1249	1264	1269	rBV	472202	1224901	27.80%	2.046%
29	9.403	1269	1272	1282	rVB2	115924	213250	4.84%	0.356%
30	9.532	1289	1294	1300	rVB2	76191	145295	3.30%	0.243%
31	9.626	1300	1310	1318	rBV3	82348	212785	4.83%	0.355%
32	9.773	1328	1335	1346	rVB3	206030	440135	9.99%	0.735%
33	9.867	1346	1351	1353	rBV	84242	128684	2.92%	0.215%
34	9.914	1353	1359	1364	rVB3	196703	350161	7.95%	0.585%
35	9.973	1364	1369	1383	rVB2	424030	1074409	24.38%	1.794%
36	10.108	1383	1392	1404	rBV	315512	767047	17.41%	1.281%

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37	10.214	1404	1410	1414	rBV5	53785	110509	2.51%	0.185%
38	10.261	1414	1418	1423	rVB4	92516	153205	3.48%	0.256%
39	10.338	1423	1431	1439	rBV	2209904	4269444	96.90%	7.130%
40	10.520	1454	1462	1468	rVB2	630895	1166711	26.48%	1.948%
41	10.673	1484	1488	1496	rBV2	97944	181257	4.11%	0.303%
42	10.761	1496	1503	1512	rBV6	167954	432648	9.82%	0.723%
43	10.855	1514	1519	1525	rVB5	86137	180837	4.10%	0.302%
44	10.950	1525	1535	1540	rBV	188708	364539	8.27%	0.609%
45	11.026	1540	1548	1549	rBV3	135023	225983	5.13%	0.377%
46	11.114	1559	1563	1566	rBV4	56951	89002	2.02%	0.149%
47	11.150	1566	1569	1572	rVV4	48464	70906	1.61%	0.118%
48	11.185	1572	1575	1579	rVV2	91974	140489	3.19%	0.235%
49	11.238	1579	1584	1589	rVV	169048	292468	6.64%	0.488%
50	11.349	1598	1603	1607	rBV2	120671	187522	4.26%	0.313%
51	11.420	1607	1615	1627	rVB4	333645	869289	19.73%	1.452%
52	11.520	1627	1632	1642	rBV	288007	527324	11.97%	0.881%
53	11.638	1645	1652	1661	rBV	1991444	3462402	78.58%	5.782%
54	11.738	1663	1669	1678	rVV	1064510	2012130	45.67%	3.360%
55	11.844	1678	1687	1698	rVV2	1219433	2824146	64.10%	4.716%
56	11.926	1698	1701	1705	rVV4	74447	106783	2.42%	0.178%
57	12.079	1723	1727	1732	rVB6	35348	62163	1.41%	0.104%
58	12.173	1732	1743	1751	rBV3	255089	788169	17.89%	1.316%
59	12.261	1753	1758	1762	rVB	141341	224515	5.10%	0.375%
60	12.344	1762	1772	1774	rBV7	84708	188397	4.28%	0.315%
61	12.385	1775	1779	1784	rVB5	84683	169299	3.84%	0.283%
62	12.473	1789	1794	1797	rBV	96913	160752	3.65%	0.268%
63	12.620	1814	1819	1829	rVB	2052028	3667096	83.23%	6.124%
64	12.714	1831	1835	1839	rBV4	49629	89937	2.04%	0.150%
65	12.808	1844	1851	1856	rVB	284852	548214	12.44%	0.916%
66	12.873	1857	1862	1866	rBV	164195	309002	7.01%	0.516%
67	12.949	1870	1875	1881	rVB2	284246	527584	11.97%	0.881%
68	13.114	1894	1903	1908	rBV2	150324	394913	8.96%	0.660%
69	13.173	1910	1913	1921	rVV4	127910	280805	6.37%	0.469%
70	13.255	1922	1927	1933	rVV	639116	1040067	23.61%	1.737%
71	13.391	1941	1950	1952	rBV7	84631	227128	5.15%	0.379%
72	13.491	1964	1967	1970	rBV3	55695	74621	1.69%	0.125%
73	13.561	1974	1979	1999	rVB2	2016071	4406082	100.00%	7.358%
74	13.726	2002	2007	2009	rBV	151051	229940	5.22%	0.384%
75	13.767	2009	2014	2019	rBV	577308	910927	20.67%	1.521%
76	13.879	2030	2033	2039	rVB4	92289	147689	3.35%	0.247%

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77	13.955	2041	2046	2050	rVV2	84439	150785	3.42%	0.252%
78	14.020	2053	2057	2065	rVV3	85713	198818	4.51%	0.332%
79	14.102	2065	2071	2077	rVV	435401	702448	15.94%	1.173%
80	14.167	2079	2082	2085	rVV	60513	74472	1.69%	0.124%
81	14.214	2085	2090	2096	rVB2	304958	522029	11.85%	0.872%
82	14.273	2096	2100	2106	rBV7	39107	79800	1.81%	0.133%
83	14.355	2109	2114	2118	rBV8	39908	72779	1.65%	0.122%
84	14.426	2118	2126	2130	rVV	225601	428834	9.73%	0.716%
85	14.467	2130	2133	2137	rVV	125785	164632	3.74%	0.275%
86	14.508	2137	2140	2144	rVB2	98311	120006	2.72%	0.200%
87	14.696	2167	2172	2177	rBV	201612	366142	8.31%	0.611%
88	14.814	2184	2192	2198	rBV3	452921	980377	22.25%	1.637%
89	14.867	2198	2201	2209	rVB2	67465	128962	2.93%	0.215%
90	14.967	2215	2218	2227	rVB8	34203	63184	1.43%	0.106%
91	15.079	2232	2237	2241	rBV3	126294	206132	4.68%	0.344%
92	15.190	2251	2256	2265	rVB2	160003	354426	8.04%	0.592%
93	15.385	2279	2289	2296	rVB	293748	601431	13.65%	1.004%
94	15.490	2302	2307	2311	rBV5	47981	88616	2.01%	0.148%
95	15.549	2313	2317	2333	rVB8	61211	216630	4.92%	0.362%
96	15.679	2333	2339	2348	rBV3	93765	209298	4.75%	0.350%
97	15.849	2361	2368	2380	rVB4	42202	138798	3.15%	0.232%
98	16.055	2397	2403	2410	rVB6	32954	80427	1.83%	0.134%
99	16.479	2469	2475	2481	rVB	82432	162862	3.70%	0.272%
100	16.685	2502	2510	2518	rBV3	44862	113014	2.56%	0.189%

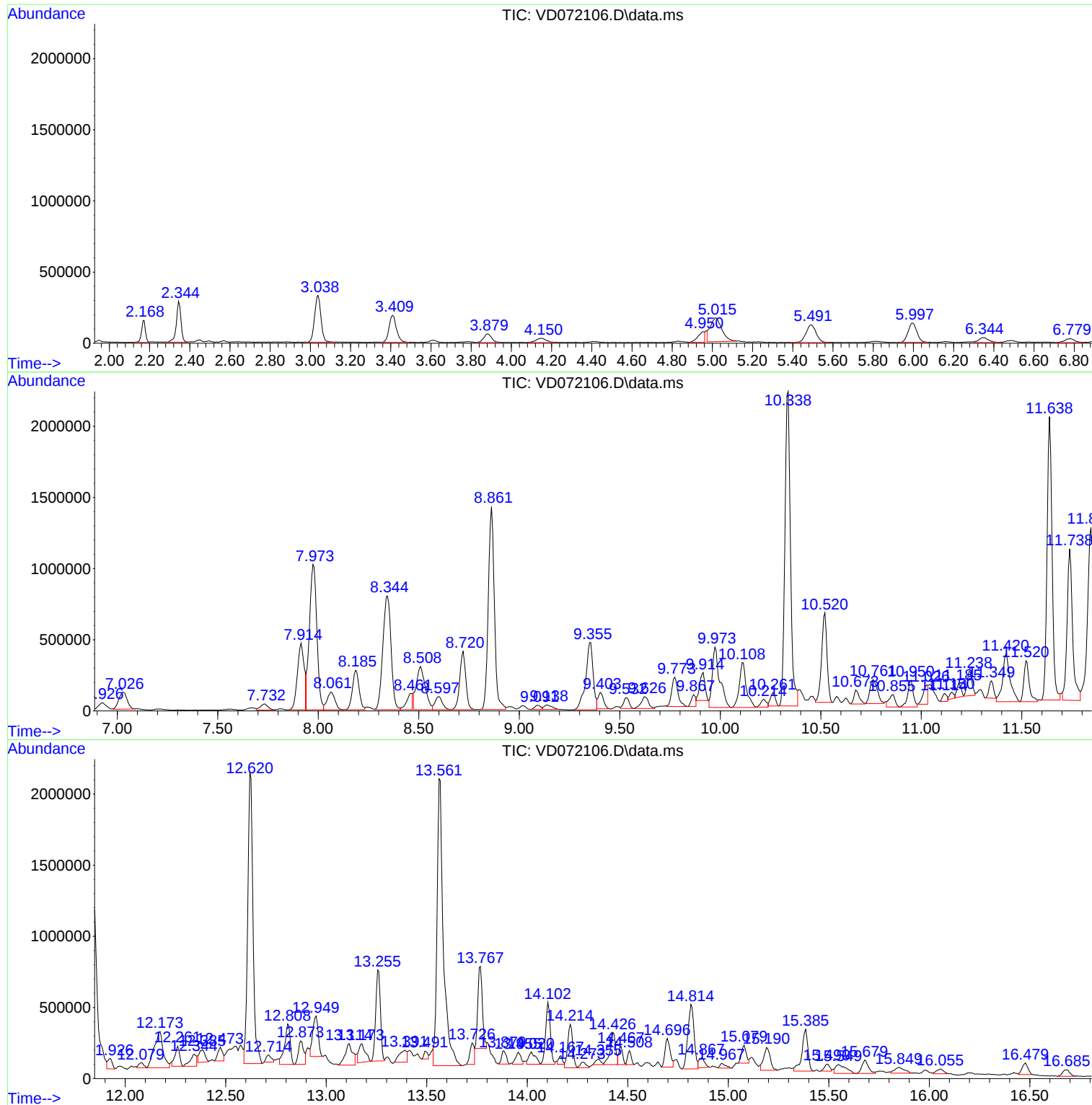
Sum of corrected areas: 59879546

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021822S.M
Quant Title : SW846 8260

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



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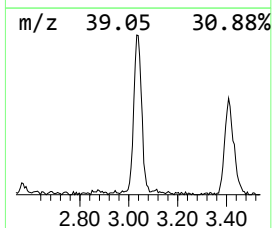
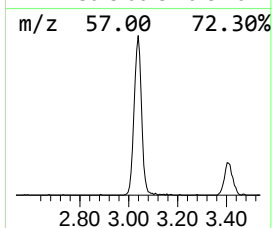
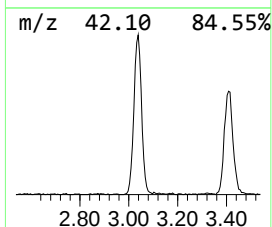
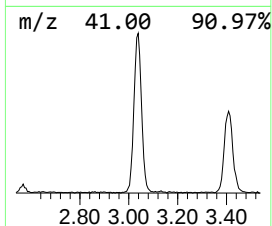
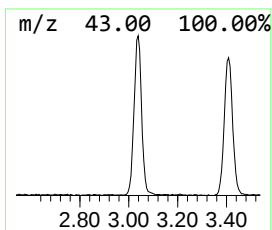
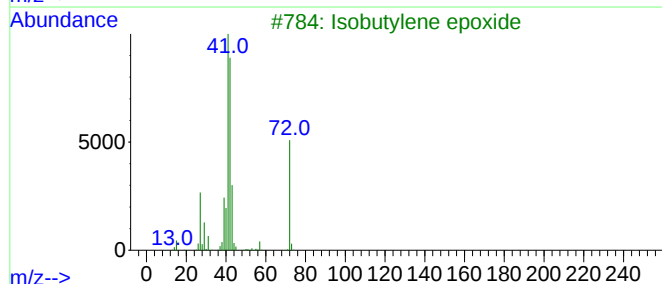
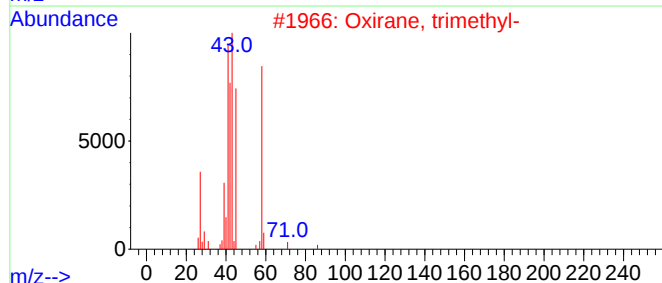
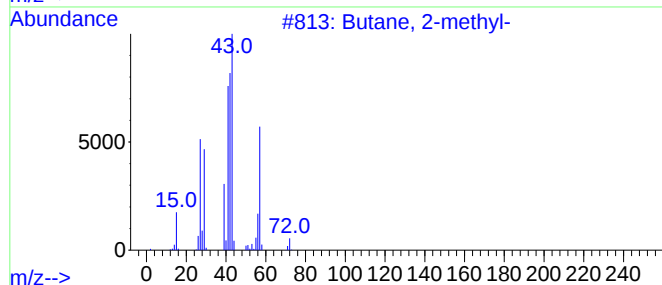
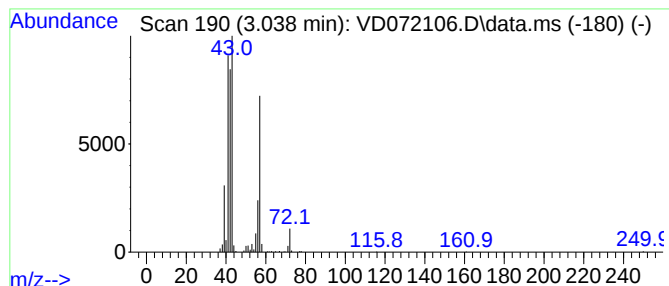
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021822S.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.038	13.66 ug/l	717699	Pentafluorobenzene	7.979

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	90
2			Oxirane, trimethyl-	86	C5H10O	005076-19-7	42
3			Isobutylene epoxide	72	C4H8O	000558-30-5	40
4			3-Buten-1-ol	72	C4H8O	000627-27-0	38
5			Pentane	72	C5H12	000109-66-0	37



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ClientSampleId :
PREM-REAR

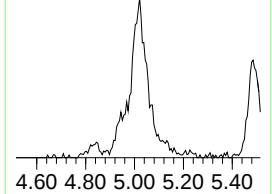
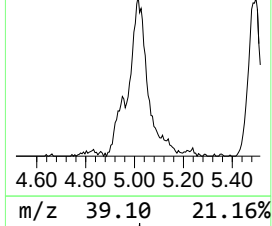
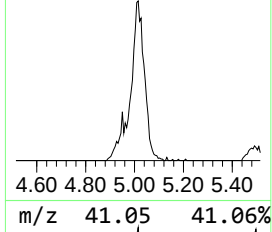
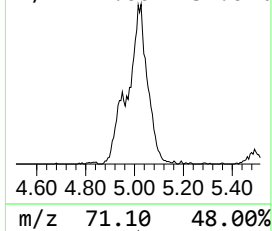
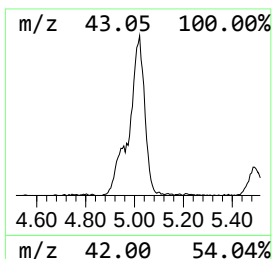
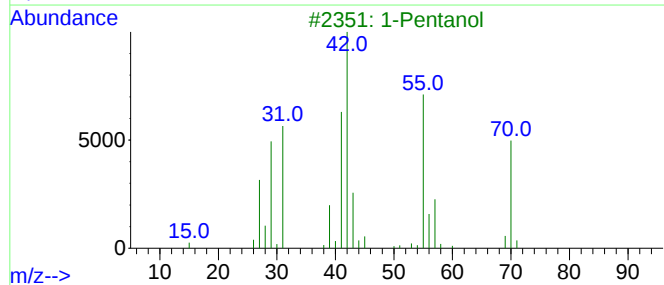
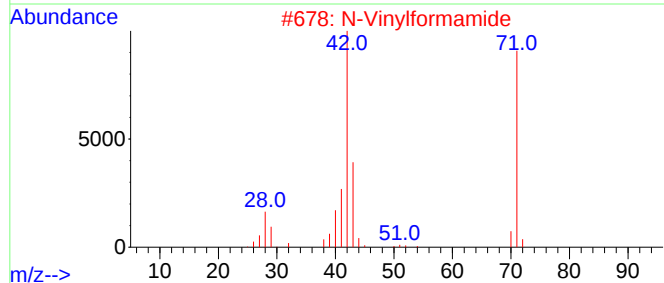
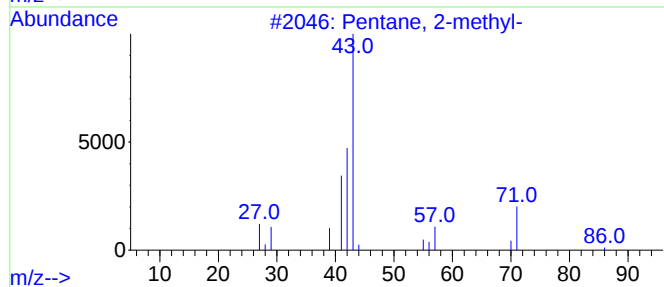
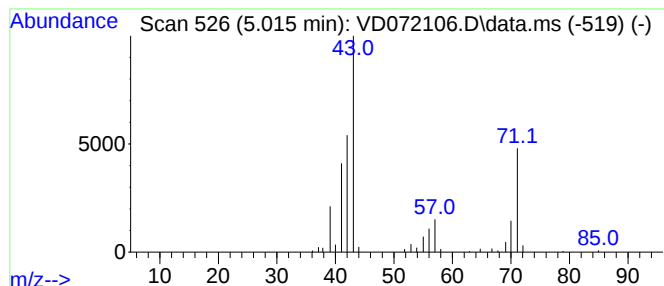
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021822S.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.015	12.50 ug/l	656847	Pentafluorobenzene	7.979

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	64
2			N-Vinylformamide	71	C3H5NO	013162-05-5	47
3			1-Pentanol	88	C5H12O	000071-41-0	38
4			Hydroxylamine, O-(3-methylbutyl)-	103	C5H13NO	019411-65-5	36
5			Butane, 2-methyl-	72	C5H12	000078-78-4	35



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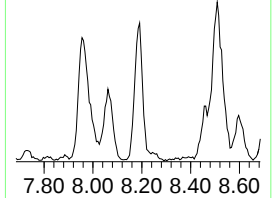
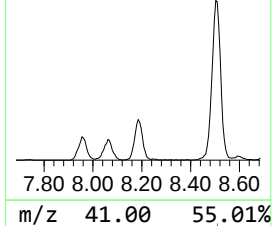
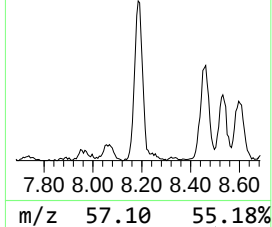
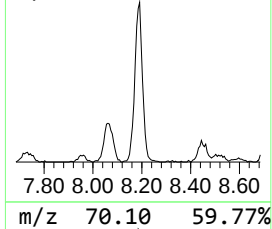
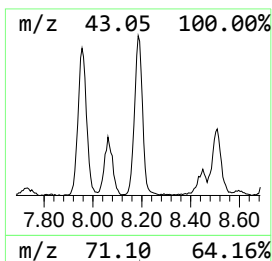
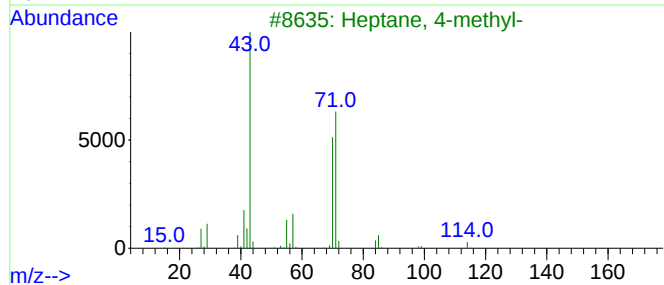
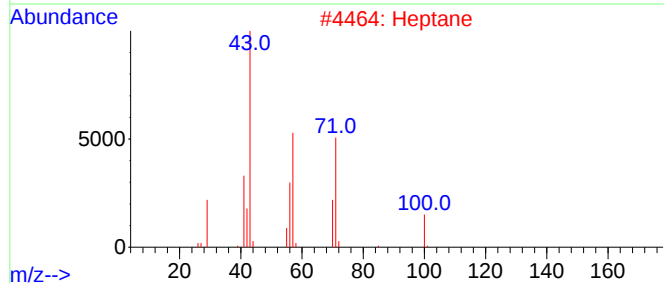
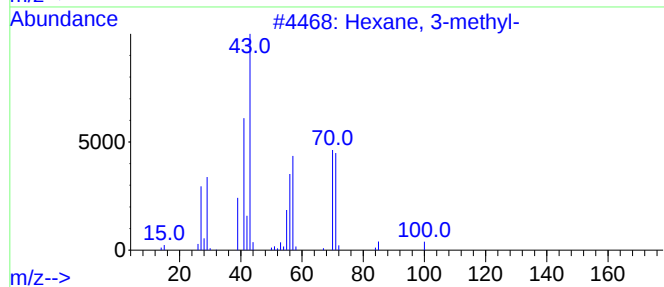
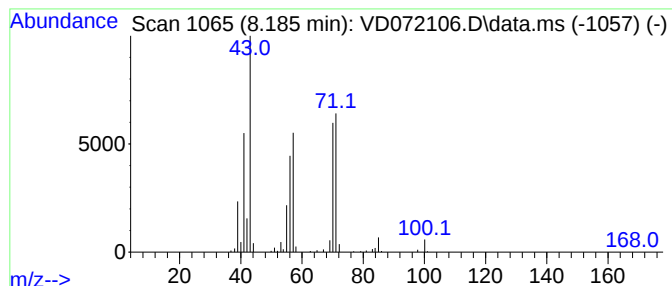
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021822S.M
Quant Title : SW846 8260

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

Peak Number 3 Hexane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.185	12.15 ug/l	638443	Pentafluorobenzene	7.979

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2			Heptane	100	C7H16	000142-82-5	80
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	53
5			1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022522\
Data File : VD072106.D
Acq On : 25 Feb 2022 17:58
Operator : VA/SY
Sample : N1655-06
Misc : 4.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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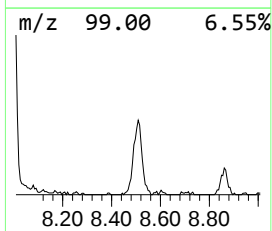
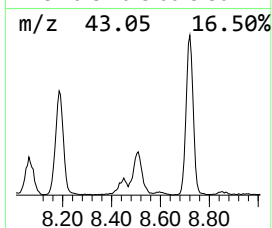
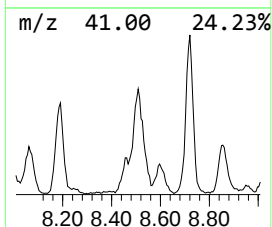
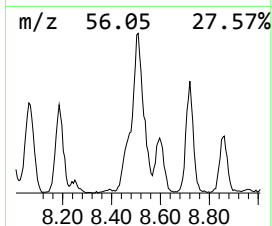
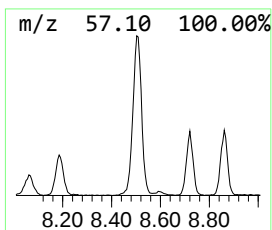
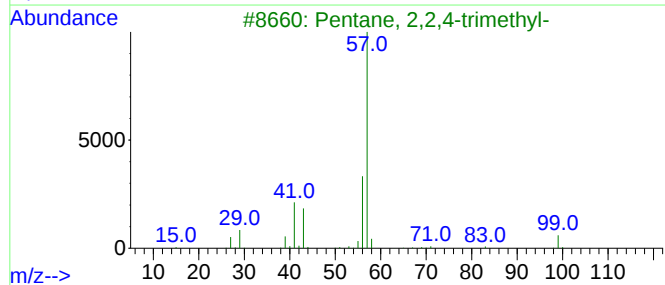
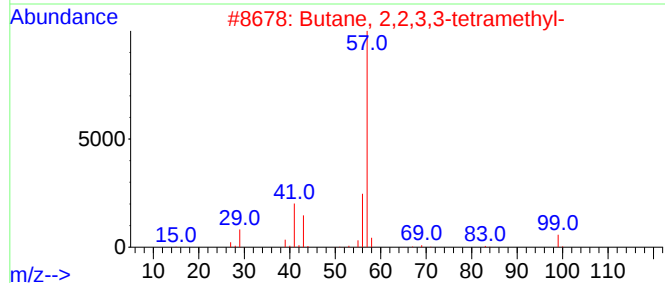
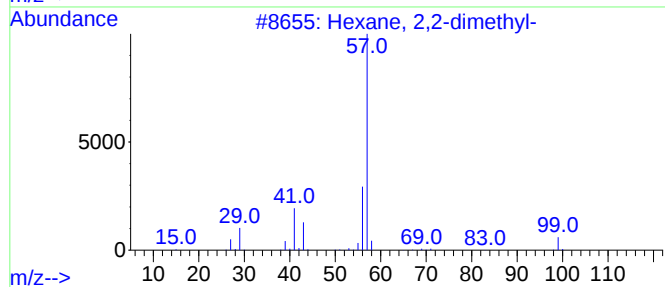
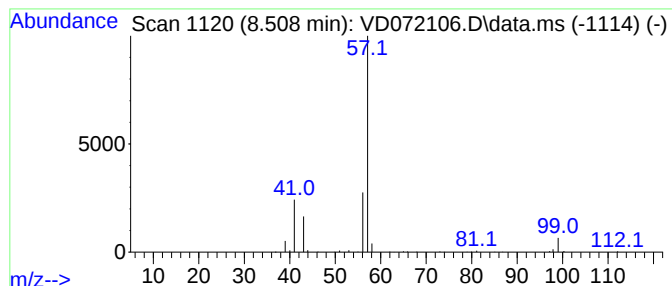
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Quant Title : SW846 8260

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

Peak Number 4 Hexane, 2,2-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.508	15.33 ug/l	914958	1,4-Difluorobenzene	8.861

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
2			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	64
3			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	50
4			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	36
5			Oxalic acid, isobutyl heptyl ester	244	C13H24O4	1000309-37-2	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022522\
Data File : VD072106.D
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Operator : VA/SY
Sample : N1655-06
Misc : 4.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PREM-REAR

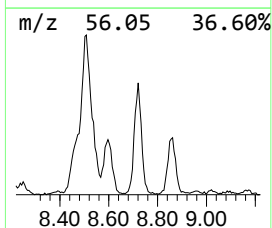
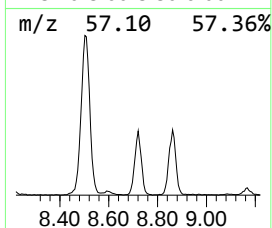
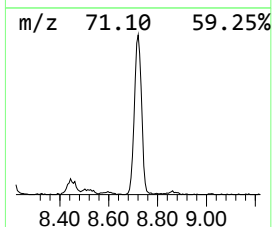
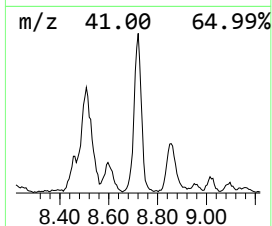
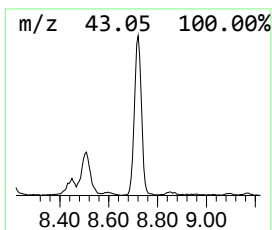
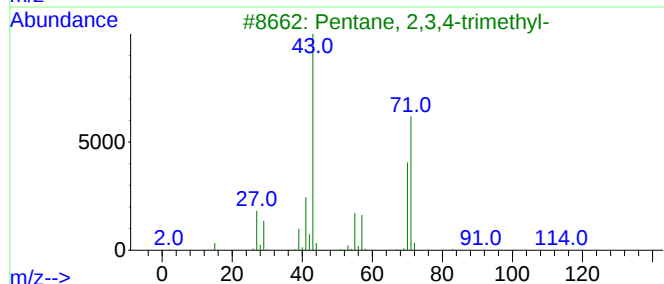
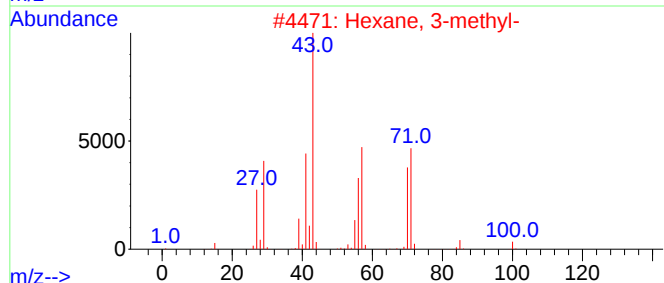
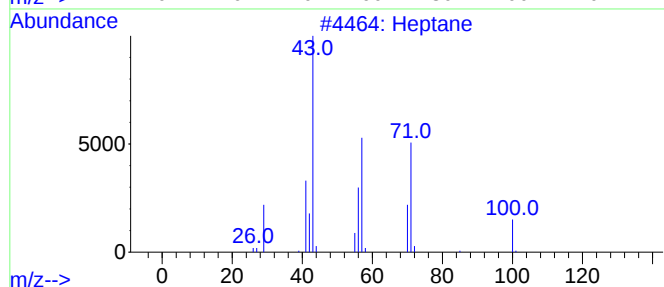
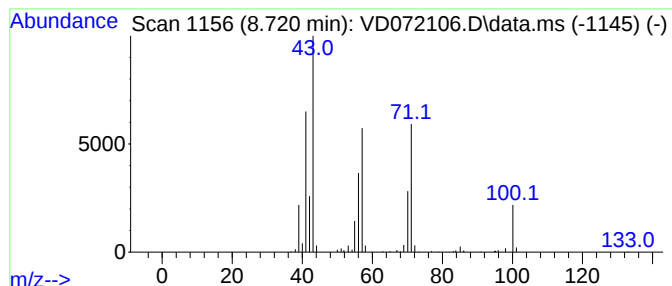
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Quant Title : SW846 8260

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

Peak Number 5 Heptane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.720	14.36 ug/l	856659	1,4-Difluorobenzene	8.861

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	91
2			Hexane, 3-methyl-	100	C7H16	000589-34-4	47
3			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	38
4			3-Hexanone	100	C6H12O	000589-38-8	38
5			Pentane, 2-methyl-	86	C6H14	000107-83-5	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022522\
Data File : VD072106.D
Acq On : 25 Feb 2022 17:58
Operator : VA/SY
Sample : N1655-06
Misc : 4.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PREM-REAR

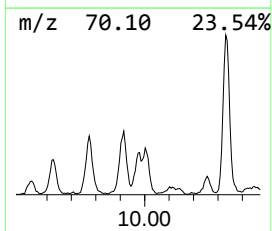
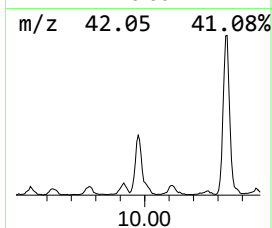
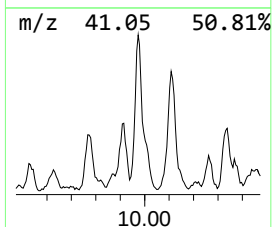
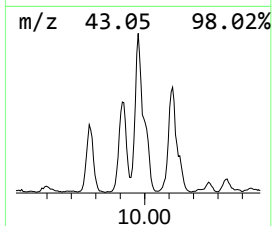
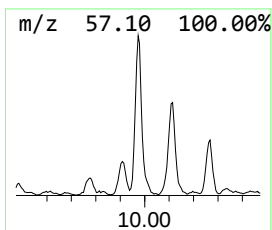
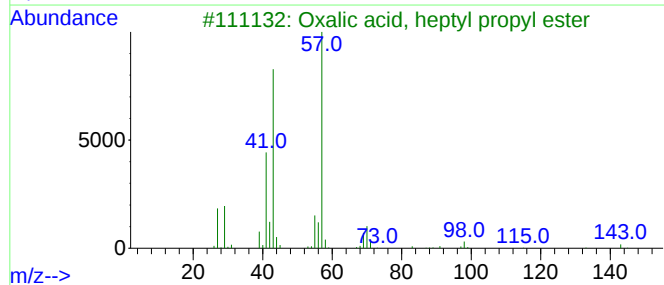
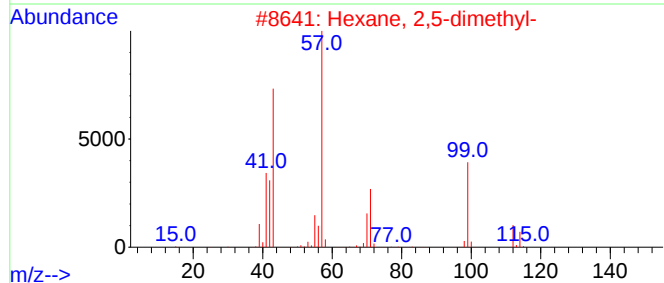
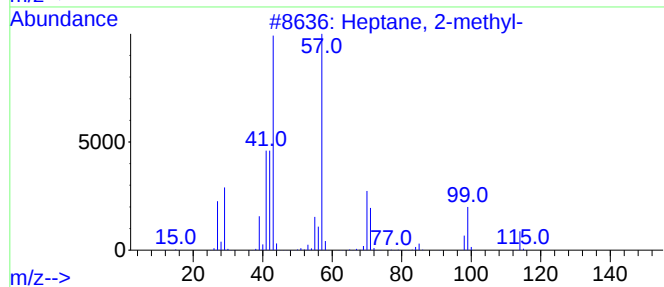
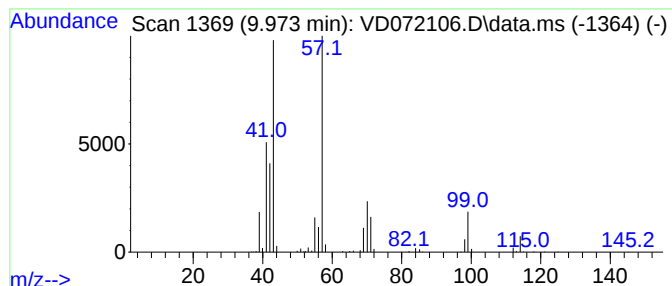
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Quant Title : SW846 8260

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

Peak Number 6 Heptane, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.973	18.01 ug/l	1074410	1,4-Difluorobenzene	8.861

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 2-methyl-	114	C8H18	000592-27-8	94
2			Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	72
3			Oxalic acid, heptyl propyl ester	230	C12H22O4	1000309-26-0	38
4			Sulfurous acid, 2-propyl heptyl ...	222	C10H22O3S	1000309-11-8	37
5			Hexane, 2,4,4-trimethyl-	128	C9H20	016747-30-1	37



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022522\
Data File : VD072106.D
Acq On : 25 Feb 2022 17:58
Operator : VA/SY
Sample : N1655-06
Misc : 4.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PREM-REAR

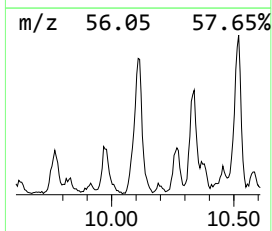
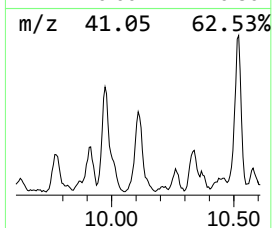
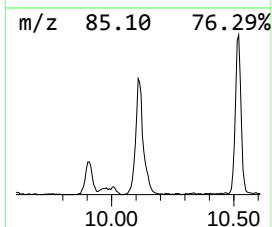
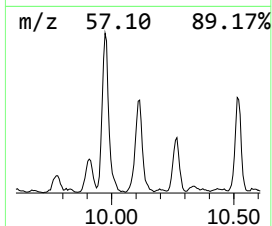
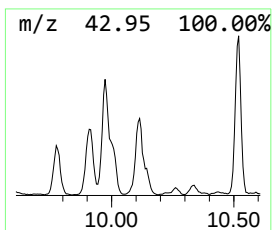
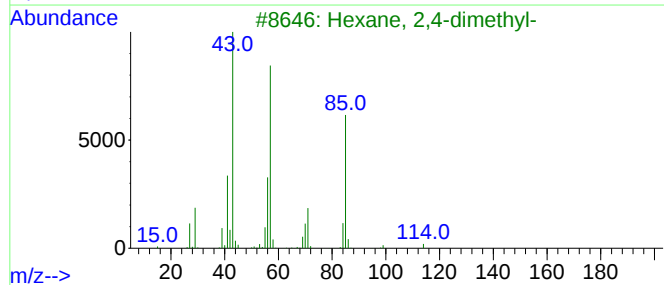
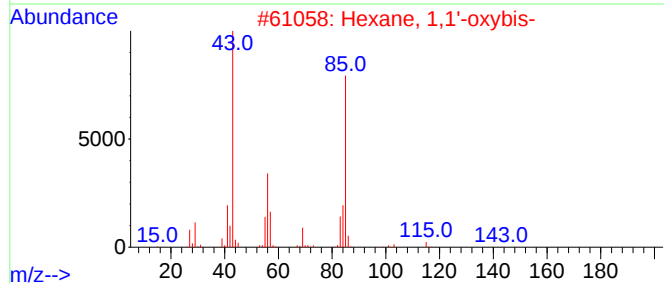
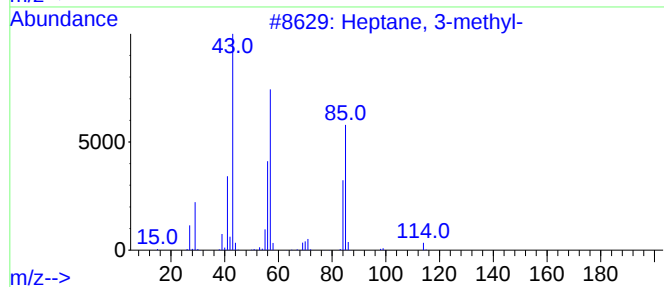
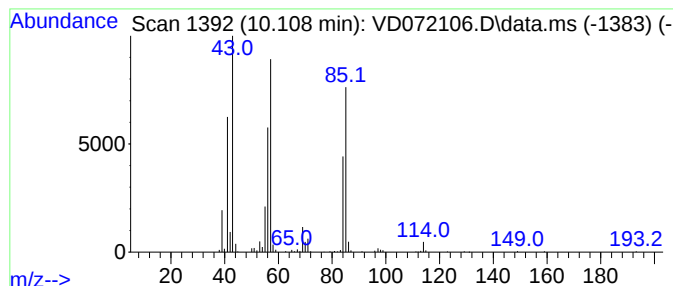
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Quant Title : SW846 8260

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

Peak Number 7 Heptane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.108	12.85 ug/l	767047	1,4-Difluorobenzene	8.861

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-methyl-	114	C8H18	000589-81-1	91
2			Hexane, 1,1'-oxybis-	186	C12H26O	000112-58-3	59
3			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	59
4			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	53
5			Undecane, 2,4-dimethyl-	184	C13H28	017312-80-0	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022522\
Data File : VD072106.D
Acq On : 25 Feb 2022 17:58
Operator : VA/SY
Sample : N1655-06
Misc : 4.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
PREM-REAR

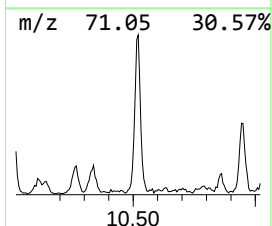
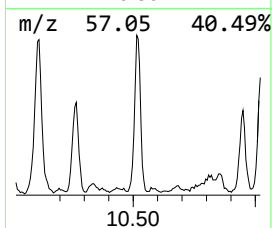
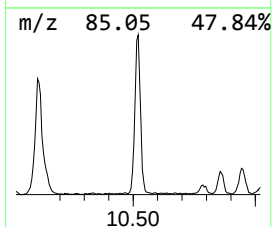
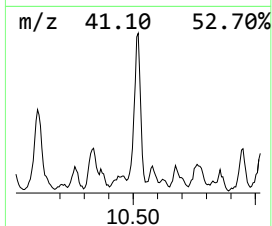
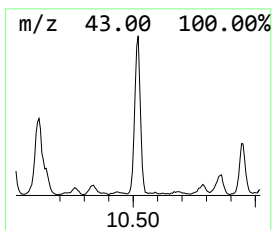
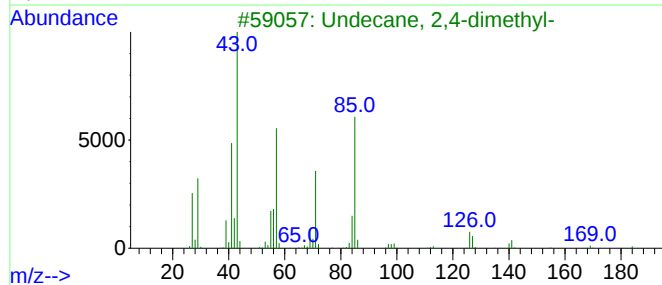
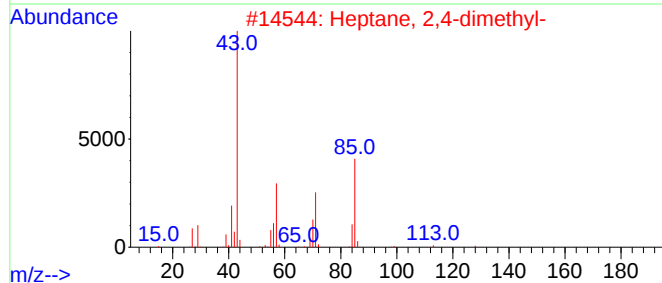
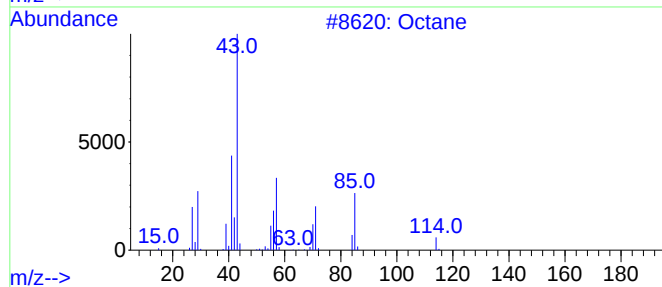
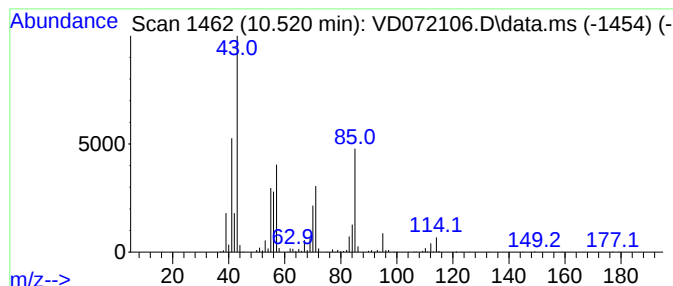
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Quant Title : SW846 8260

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

Peak Number 8 Octane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.520	16.85 ug/l	1166710	Chlorobenzene-d5	11.638

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane			114	C8H18	000111-65-9	90
2	Heptane, 2,4-dimethyl-			128	C9H20	002213-23-2	72
3	Undecane, 2,4-dimethyl-			184	C13H28	017312-80-0	53
4	Nonane			128	C9H20	000111-84-2	53
5	Hexane, 3,3-dimethyl-			114	C8H18	000563-16-6	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022522\
Data File : VD072106.D
Acq On : 25 Feb 2022 17:58
Operator : VA/SY
Sample : N1655-06
Misc : 4.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PREM-REAR

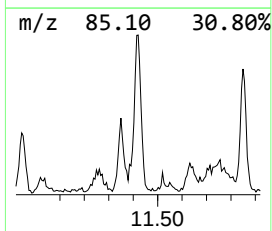
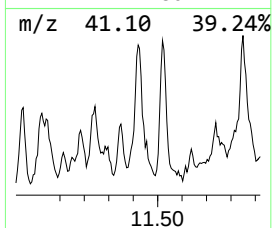
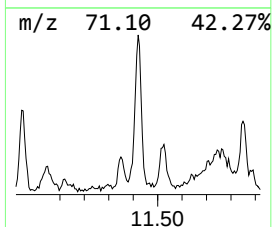
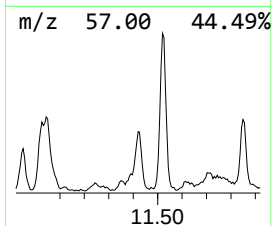
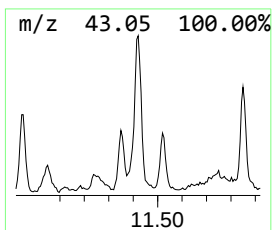
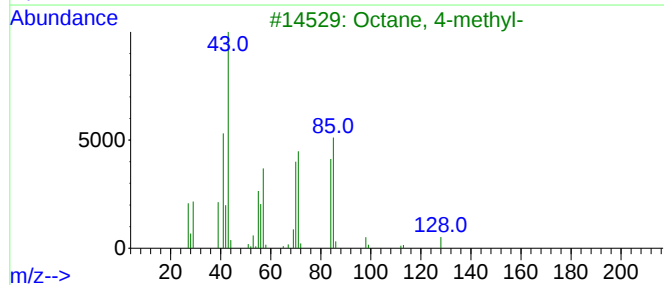
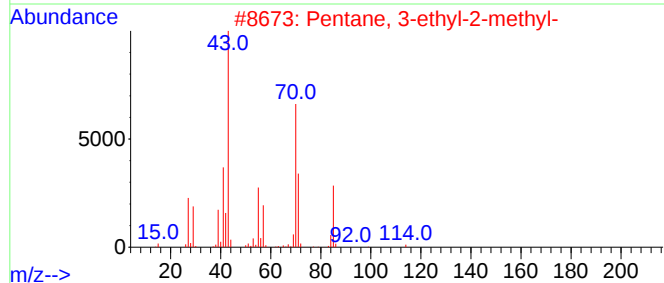
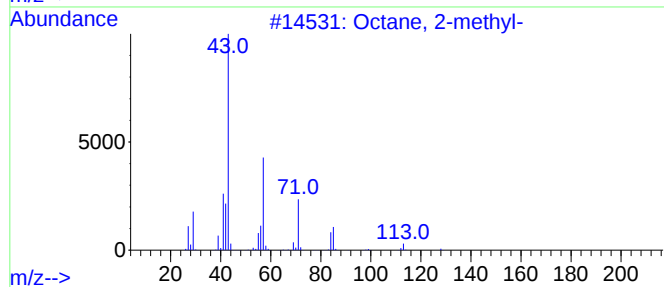
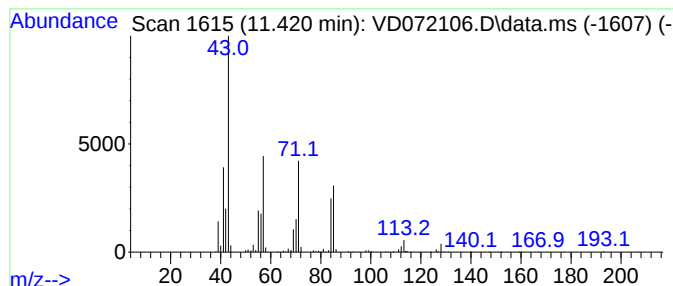
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Quant Title : SW846 8260

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

Peak Number 9 Octane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.420	12.55 ug/l	869289	Chlorobenzene-d5	11.638

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2-methyl-	128	C9H20	003221-61-2	64
2			Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	64
3			Octane, 4-methyl-	128	C9H20	002216-34-4	53
4			Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	53
5			Ether, hexyl pentyl	172	C11H24O	032357-83-8	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022522\
Data File : VD072106.D
Acq On : 25 Feb 2022 17:58
Operator : VA/SY
Sample : N1655-06
Misc : 4.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

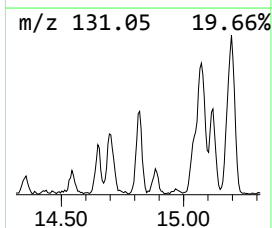
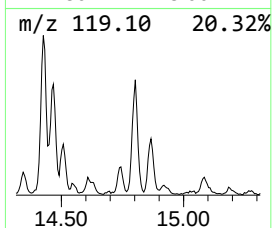
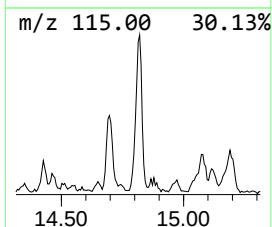
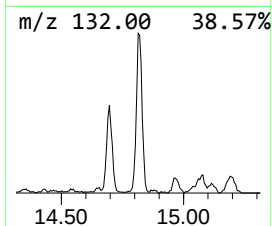
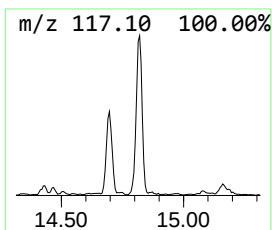
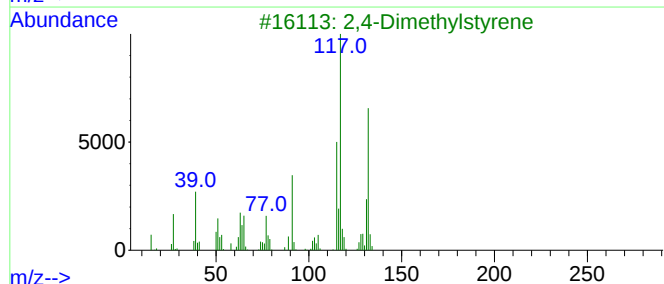
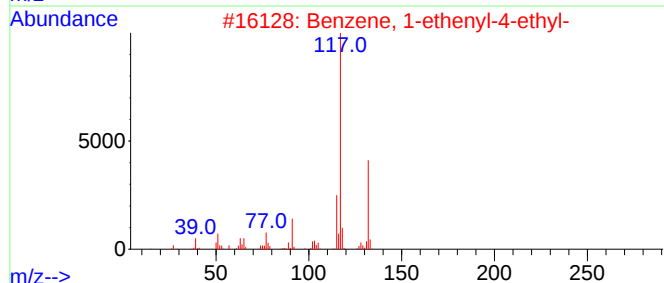
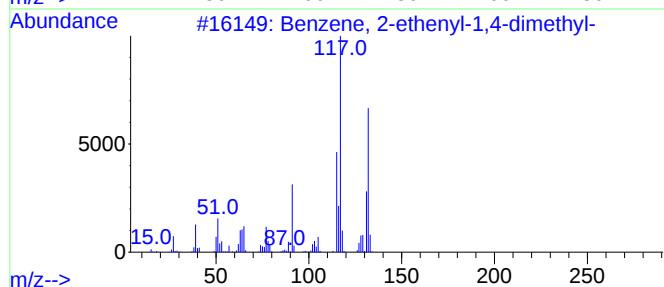
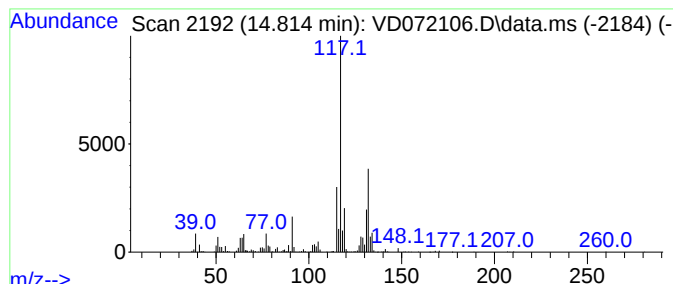
Instrument :
MSVOA_D
ClientSampleId :
PREM-REAR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021822S.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.814	11.13 ug/l	980377	1,4-Dichlorobenzene-d4	13.561	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
3	2,4-Dimethylstyrene	132	C10H12	002234-20-0	91
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022522\
Data File : VD072106.D
Acq On : 25 Feb 2022 17:58
Operator : VA/SY
Sample : N1655-06
Misc : 4.43G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PREM-REAR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021822S.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
Butane, 2-methyl-	3.038	13.7	ug/l	717699	1	7.979	2627940	50.0
Pentane, 2-methyl-	5.015	12.5	ug/l	656847	1	7.979	2627940	50.0
Hexane, 3-methyl-	8.185	12.2	ug/l	638443	1	7.979	2627940	50.0
Hexane, 2,2-dim...	8.508	15.3	ug/l	914958	2	8.861	2983580	50.0
Heptane	8.720	14.4	ug/l	856659	2	8.861	2983580	50.0
Heptane, 2-methyl-	9.973	18.0	ug/l	1074410	2	8.861	2983580	50.0
Heptane, 3-methyl-	10.108	12.9	ug/l	767047	2	8.861	2983580	50.0
Octane	10.520	16.9	ug/l	1166710	3	11.638	3462400	50.0
Octane, 2-methyl-	11.420	12.6	ug/l	869289	3	11.638	3462400	50.0
Benzene, 2-ethe...	14.814	11.1	ug/l	980377	4	13.561	4406080	50.0