

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
 Data File : VD068710.D
 Acq On : 29 Mar 2021 20:50
 Operator : VA/SY
 Sample : M1836-01
 Misc : 8.84G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 B1-0-2

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\82D031921S.M

Title : SW846 8260

Signal : TIC: VD068710.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.167	35	42	52	rBV	432012	581000	1.34%	0.138%
2	2.344	61	72	82	rBV	2966802	4678397	10.79%	1.113%
3	3.032	177	189	219	rBV	10781477	22867976	52.72%	5.440%
4	3.408	243	253	275	rVB	3798464	9511709	21.93%	2.263%
5	3.879	323	333	350	rVB	2757406	7161414	16.51%	1.703%
6	4.138	364	377	399	rVB2	454565	1536170	3.54%	0.365%
7	4.949	502	515	519	rVV2	1662694	5693660	13.13%	1.354%
8	5.014	520	526	555	rVV4	3103672	13585775	31.32%	3.232%
9	5.491	590	607	635	rBV2	2598708	9146881	21.09%	2.176%
10	5.808	649	661	677	rVB2	209815	692750	1.60%	0.165%
11	5.996	679	693	711	rBV	1963942	5936756	13.69%	1.412%
12	6.343	731	752	764	rBV	1266600	3931602	9.06%	0.935%
13	6.485	764	776	786	rBV2	558514	1703993	3.93%	0.405%
14	6.779	814	826	839	rBV2	1168018	3438597	7.93%	0.818%
15	6.926	839	851	859	rBV	1551881	4635121	10.69%	1.103%
16	7.026	859	868	877	rBV	3037097	8471884	19.53%	2.015%
17	7.732	981	988	997	rVB	1054302	2575586	5.94%	0.613%
18	7.967	1008	1028	1038	rBV6	3204513	11834018	27.28%	2.815%
19	8.067	1038	1045	1056	rVB	3402678	8463346	19.51%	2.013%
20	8.190	1056	1066	1074	rBV	3079407	7160353	16.51%	1.703%
21	8.355	1083	1094	1104	rBV	13954448	32219245	74.28%	7.664%
22	8.508	1104	1120	1131	rBV	16249243	43375424	100.00%	10.318%
23	8.602	1131	1136	1146	rVB3	820541	1946857	4.49%	0.463%
24	8.720	1146	1156	1170	rVB2	2091036	4624065	10.66%	1.100%
25	8.861	1170	1180	1186	rBV2	2821057	6743090	15.55%	1.604%
26	9.020	1202	1207	1214	rVB	427831	800636	1.85%	0.190%
27	9.096	1214	1220	1223	rBV2	336684	667298	1.54%	0.159%
28	9.137	1223	1227	1249	rVB2	609237	1508851	3.48%	0.359%
29	9.355	1249	1264	1269	rBV	5170133	12359414	28.49%	2.940%
30	9.408	1269	1273	1280	rVB	2434679	4567074	10.53%	1.086%
31	9.490	1280	1287	1291	rBV	992249	1991686	4.59%	0.474%
32	9.537	1291	1295	1301	rVB	777029	1382546	3.19%	0.329%
33	9.631	1301	1311	1317	rVB3	375374	985031	2.27%	0.234%
34	9.784	1328	1337	1346	rVB	10898256	21867592	50.41%	5.202%
35	9.914	1346	1359	1366	rBV2	14797678	32068484	73.93%	7.628%
36	9.979	1366	1370	1374	rBV	1095421	1640166	3.78%	0.390%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
 Data File : VD068710.D
 Acq On : 29 Mar 2021 20:50
 Operator : VA/SY
 Sample : M1836-01
 Misc : 8.84G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 B1-0-2

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\82D031921S.M
 Title : SW846 8260

37	10.114	1383	1393	1405	rBV	2028130	5139814	11.85%	1.223%
38	10.220	1405	1411	1414	rBV	577722	1003677	2.31%	0.239%
39	10.267	1414	1419	1425	rVB2	1848730	3336484	7.69%	0.794%
40	10.337	1425	1431	1440	rBV	2511511	5018525	11.57%	1.194%
41	10.526	1455	1463	1469	rVB5	1335319	2962474	6.83%	0.705%
42	10.678	1485	1489	1497	rBV3	314692	749591	1.73%	0.178%
43	10.761	1497	1503	1515	rVB5	862067	2532501	5.84%	0.602%
44	10.861	1515	1520	1526	rVB2	328813	642529	1.48%	0.153%
45	10.949	1526	1535	1541	rBV3	263092	650969	1.50%	0.155%
46	11.031	1541	1549	1560	rBV3	717894	2404740	5.54%	0.572%
47	11.155	1566	1570	1573	rVV2	324060	547453	1.26%	0.130%
48	11.190	1573	1576	1581	rVV	326516	566242	1.31%	0.135%
49	11.355	1599	1604	1607	rVV2	278311	433783	1.00%	0.103%
50	11.426	1607	1616	1628	rVB3	816425	2222414	5.12%	0.529%
51	11.525	1628	1633	1644	rBV	768214	1453971	3.35%	0.346%
52	11.643	1645	1653	1659	rVV	1847118	3291739	7.59%	0.783%
53	11.743	1661	1670	1679	rVV	8886223	14829335	34.19%	3.527%
54	11.855	1679	1689	1699	rVV	5239794	11097984	25.59%	2.640%
55	12.178	1733	1744	1752	rBV3	390727	1091800	2.52%	0.260%
56	12.349	1763	1773	1777	rBV4	172584	444755	1.03%	0.106%
57	12.478	1789	1795	1801	rBV	1821072	2980984	6.87%	0.709%
58	12.625	1813	1820	1831	rVB2	1877264	3681600	8.49%	0.876%
59	12.820	1846	1853	1858	rVB	3536358	5663591	13.06%	1.347%
60	12.878	1858	1863	1866	rVV	720640	1188263	2.74%	0.283%
61	12.914	1866	1869	1872	rVV	1184157	1717861	3.96%	0.409%
62	12.955	1872	1876	1883	rVB	1681251	2706578	6.24%	0.644%
63	13.120	1895	1904	1911	rBV	806351	1413845	3.26%	0.336%
64	13.267	1922	1929	1935	rVB	6744486	10666054	24.59%	2.537%
65	13.396	1942	1951	1955	rBV3	388708	1097307	2.53%	0.261%
66	13.602	1974	1986	1995	rVB2	4134429	9416229	21.71%	2.240%
67	13.731	2002	2008	2011	rBV	730571	1217949	2.81%	0.290%
68	13.772	2011	2015	2019	rVV	2555780	4468915	10.30%	1.063%
69	13.808	2019	2021	2031	rVV4	886833	1715038	3.95%	0.408%
70	13.967	2042	2048	2053	rVV	538881	838172	1.93%	0.199%
71	14.025	2053	2058	2061	rVV	509002	858670	1.98%	0.204%
72	14.055	2061	2063	2067	rVV	425598	540996	1.25%	0.129%
73	14.114	2067	2073	2079	rVV	1456374	2214759	5.11%	0.527%
74	14.225	2086	2092	2097	rVB2	900419	1511175	3.48%	0.359%
75	14.355	2109	2114	2119	rBV2	286741	442788	1.02%	0.105%
76	14.437	2119	2128	2131	rBV	515183	919093	2.12%	0.219%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
Data File : VD068710.D
Acq On : 29 Mar 2021 20:50
Operator : VA/SY
Sample : M1836-01
Misc : 8.84G/5.00ml/MSVOA_D/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B1-0-2

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\82D031921S.M
Title : SW846 8260

77	14.472	2131	2134	2139	rVB	510933	724917	1.67%	0.172%
78	14.708	2169	2174	2178	rBV	300548	501353	1.16%	0.119%
79	14.825	2186	2194	2199	rBV2	538863	1132893	2.61%	0.269%

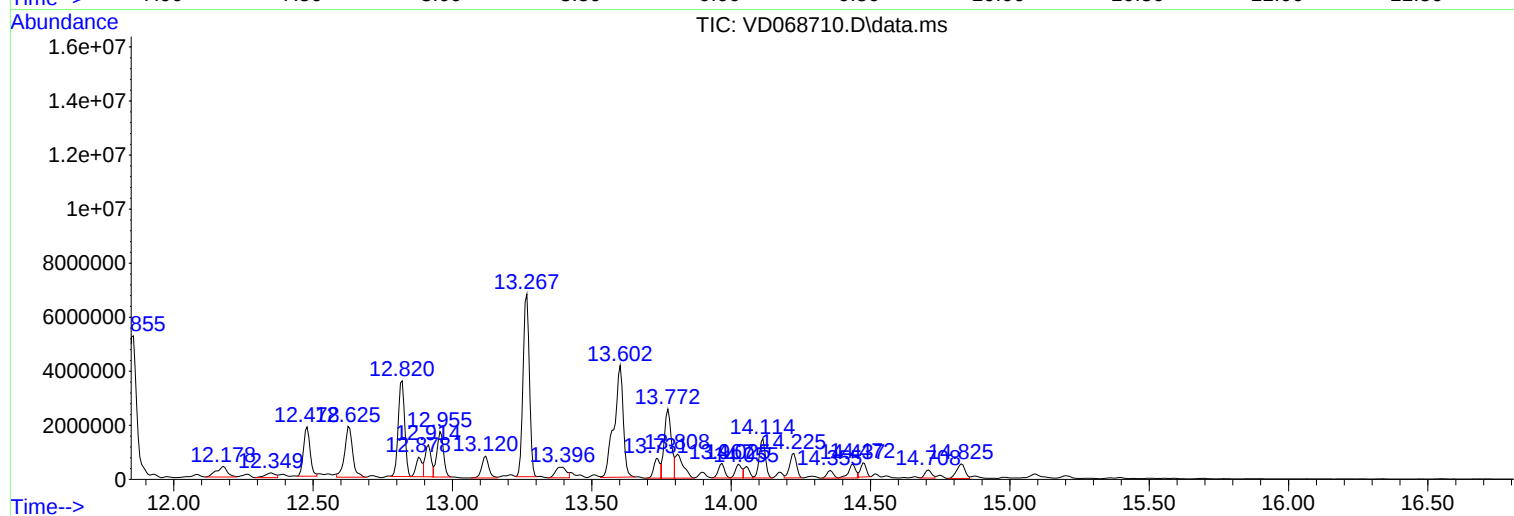
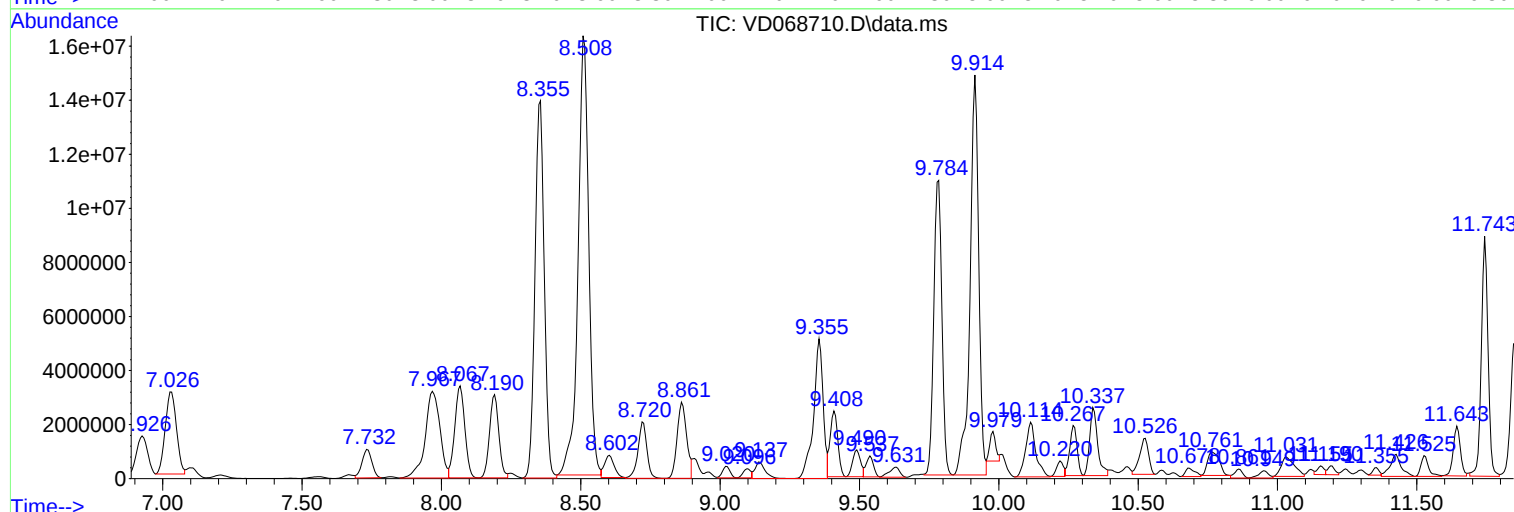
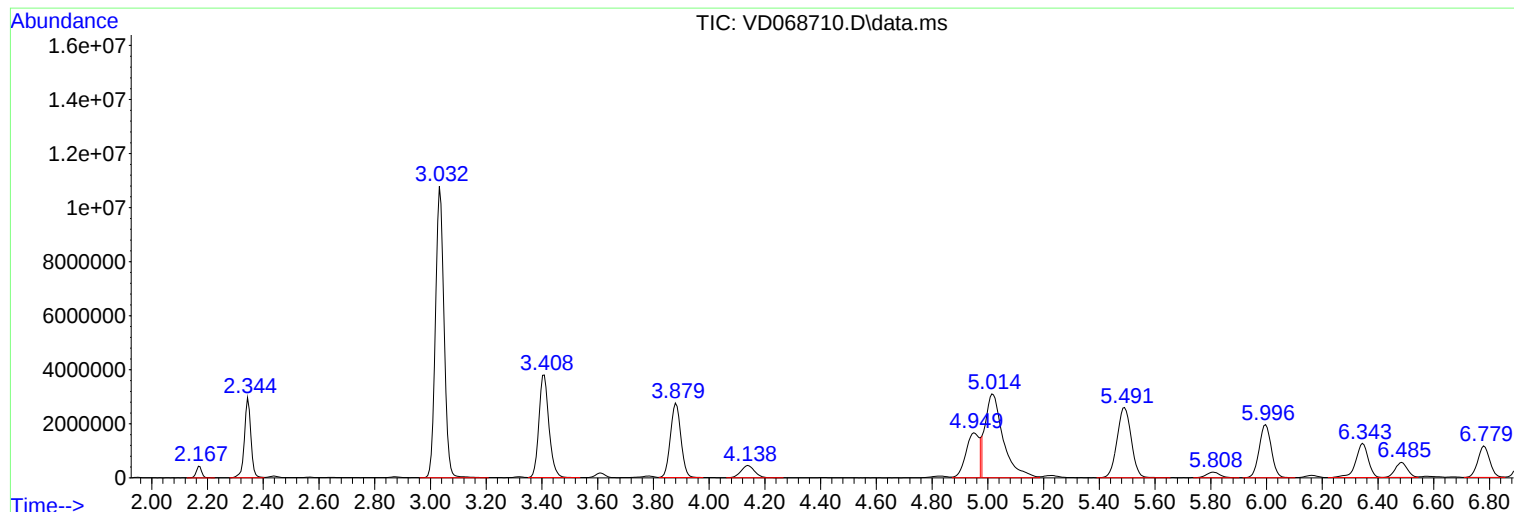
Sum of corrected areas: 420394257

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
Data File : VD068710.D
Acq On : 29 Mar 2021 20:50
Operator : VA/SY
Sample : M1836-01
Misc : 8.84G/5.00ml/MSVOA_D/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B1-0-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031921S.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
Data File : VD068710.D
Acq On : 29 Mar 2021 20:50
Operator : VA/SY
Sample : M1836-01
Misc : 8.84G/5.00ml/MSVOA_D/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B1-0-2

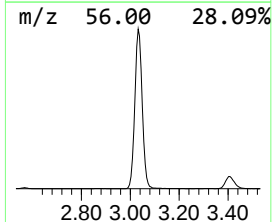
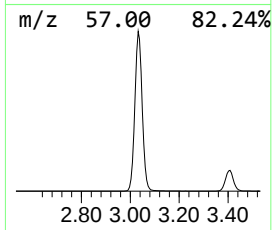
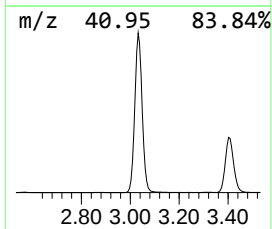
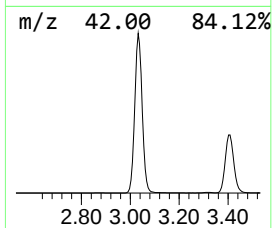
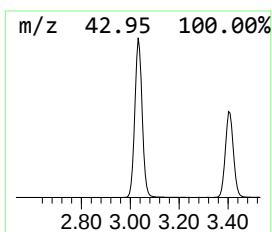
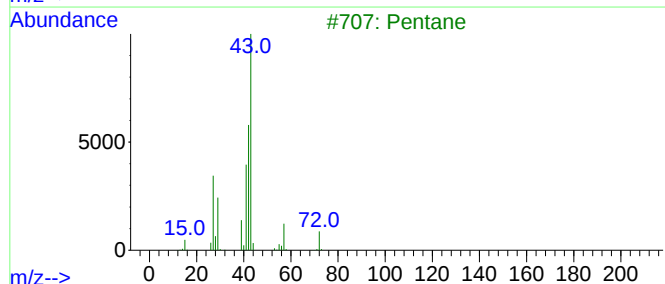
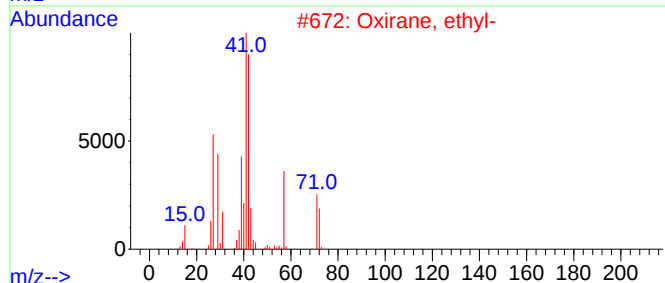
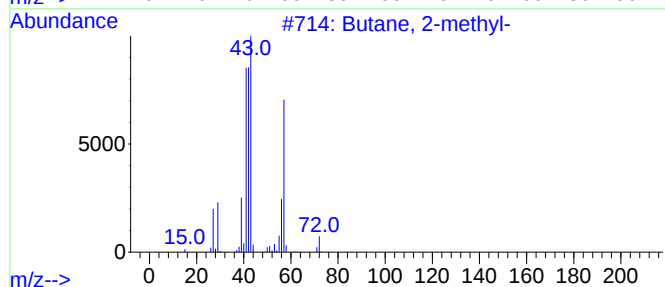
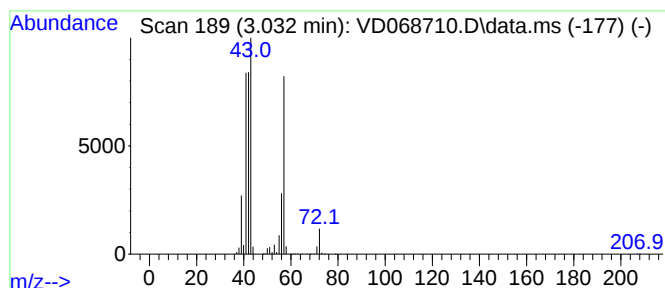
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031921S.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.032	96.62 ug/l	22868000	Pentafluorobenzene	7.979

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	94
2			Oxirane, ethyl-	72	C4H8O	000106-88-7	47
3			Pentane	72	C5H12	000109-66-0	42
4			1,3-Dimethyldiaziridine	72	C3H8N2	026177-36-6	33
5			Oxirane, trimethyl-	86	C5H10O	005076-19-7	33



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
Data File : VD068710.D
Acq On : 29 Mar 2021 20:50
Operator : VA/SY
Sample : M1836-01
Misc : 8.84G/5.00ml/MSVOA_D/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B1-0-2

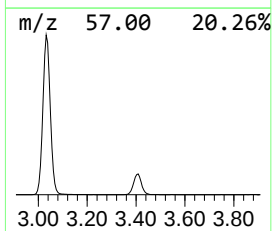
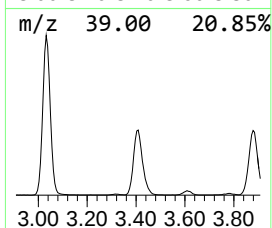
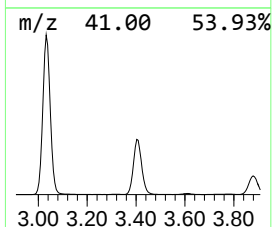
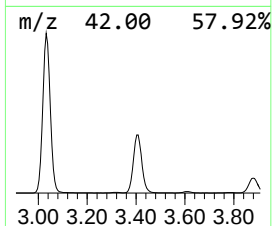
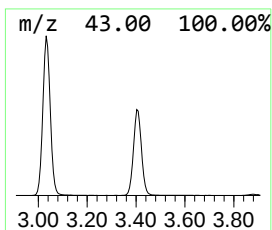
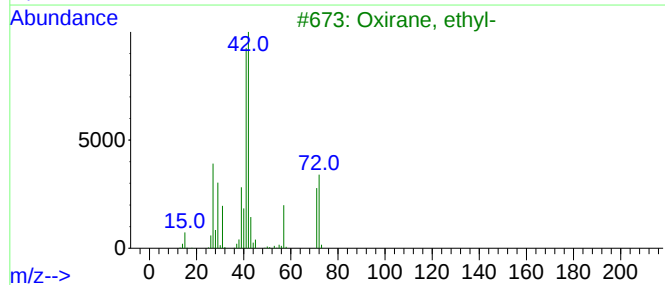
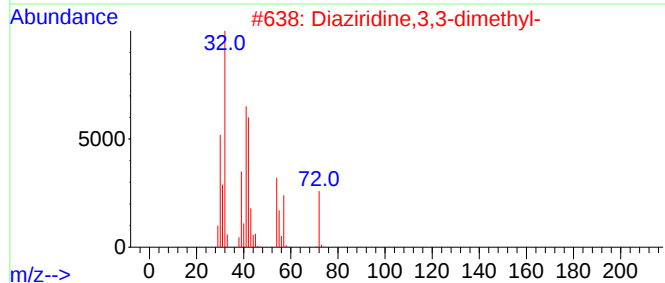
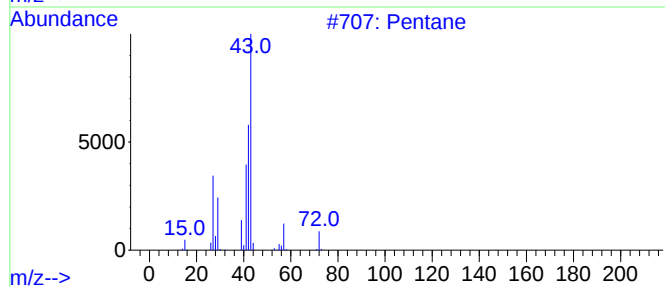
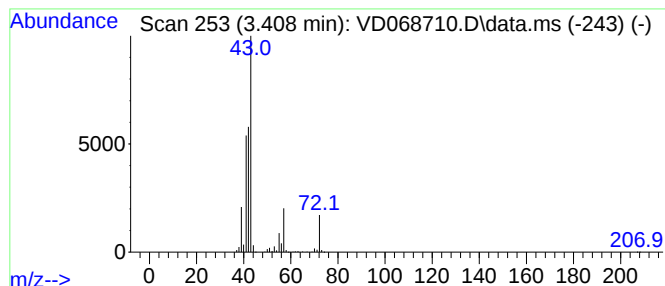
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031921S.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.408	40.19 ug/l	9511710	Pentafluorobenzene	7.979

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane	72	C5H12	000109-66-0	86
2			Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	43
3			Oxirane, ethyl-	72	C4H8O	000106-88-7	40
4			2-Butanone	72	C4H8O	000078-93-3	16
5			Butanal, 2-ethyl-	100	C6H12O	000097-96-1	10



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
Data File : VD068710.D
Acq On : 29 Mar 2021 20:50
Operator : VA/SY
Sample : M1836-01
Misc : 8.84G/5.00ml/MSVOA_D/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B1-0-2

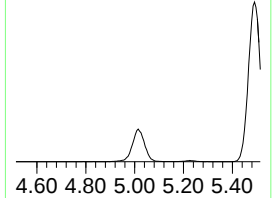
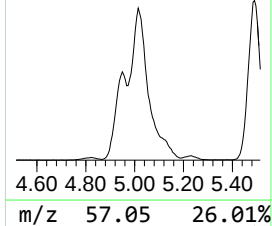
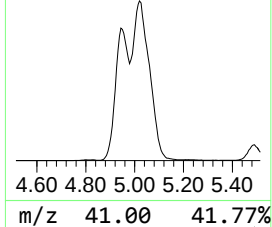
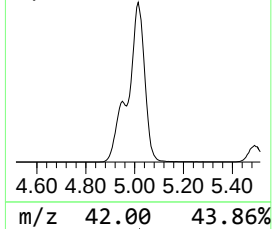
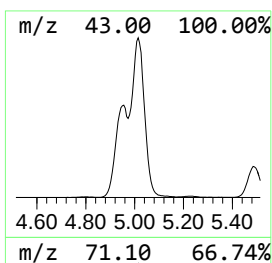
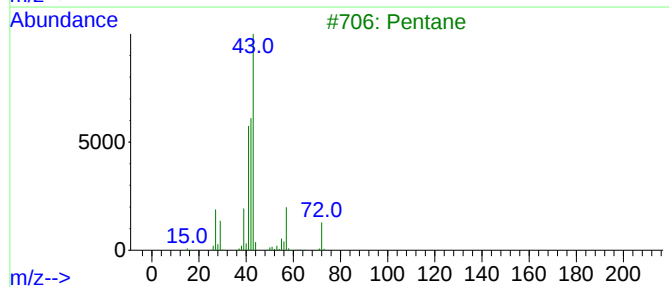
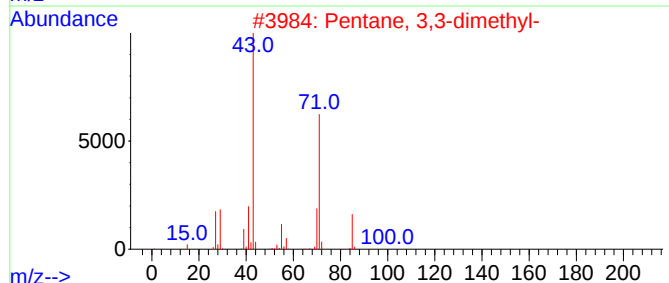
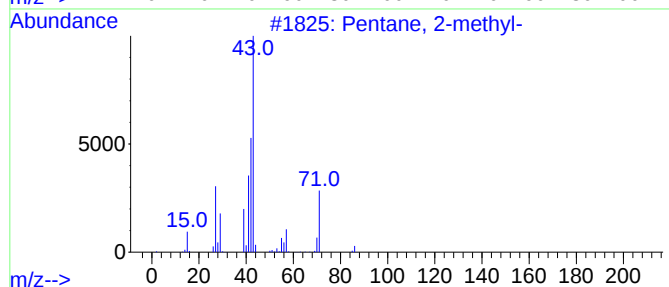
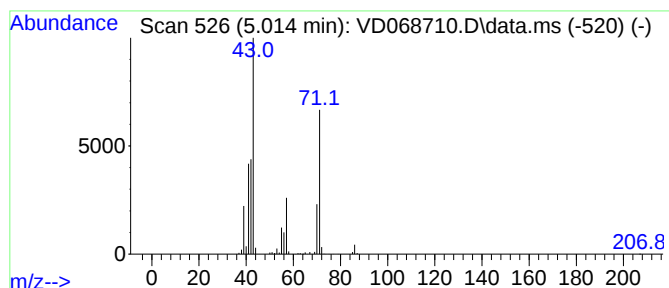
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031921S.M
Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.014	57.40 ug/l	13585800	Pentafluorobenzene	7.979

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	53
2			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	47
3			Pentane	72	C5H12	000109-66-0	38
4			Butane, 2-methyl-	72	C5H12	000078-78-4	38
5			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	33



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
Data File : VD068710.D
Acq On : 29 Mar 2021 20:50
Operator : VA/SY
Sample : M1836-01
Misc : 8.84G/5.00ml/MSVOA_D/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B1-0-2

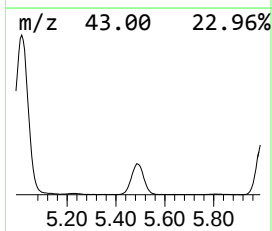
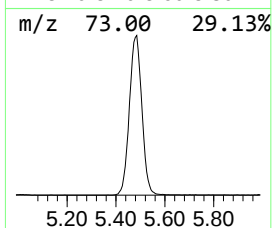
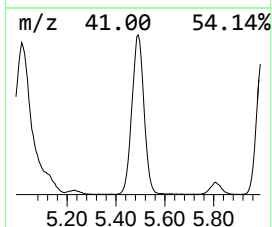
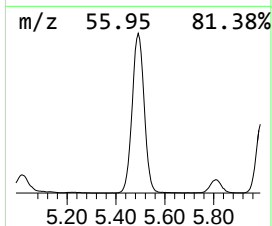
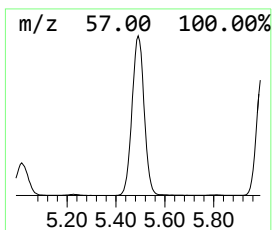
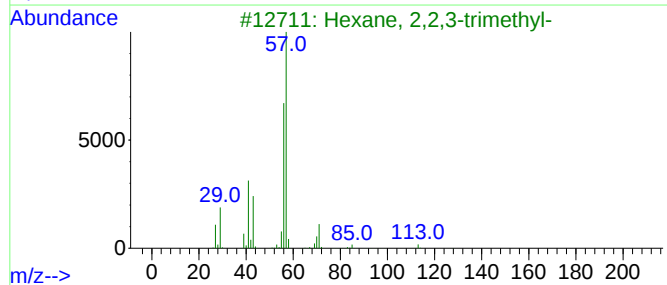
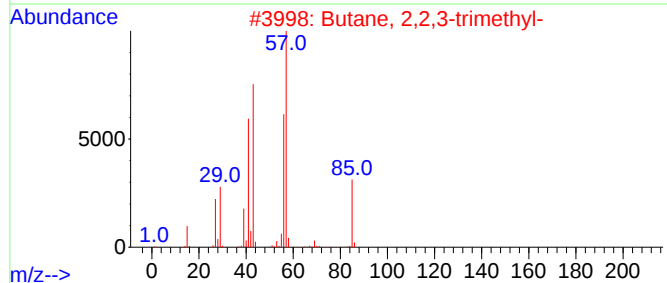
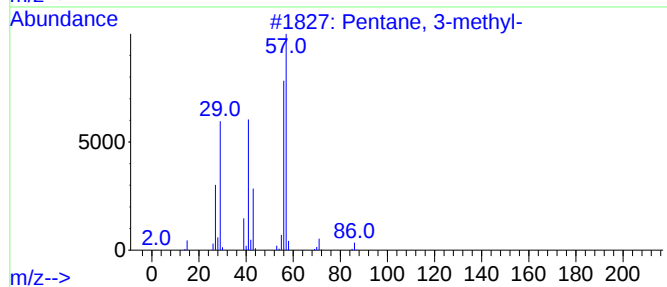
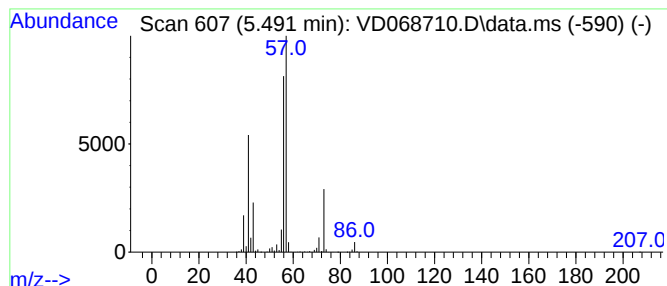
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031921S.M
Quant Title : SW846 8260

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

Peak Number 4 Pentane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.491	38.65 ug/l	9146880	Pentafluorobenzene	7.979

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	83
2			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64
3			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	50
4			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	50
5			Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
Data File : VD068710.D
Acq On : 29 Mar 2021 20:50
Operator : VA/SY
Sample : M1836-01
Misc : 8.84G/5.00ml/MSVOA_D/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B1-0-2

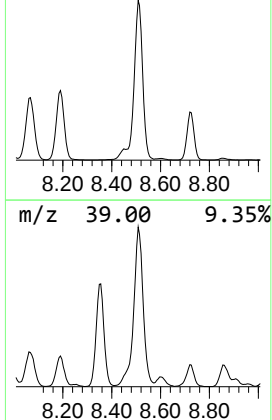
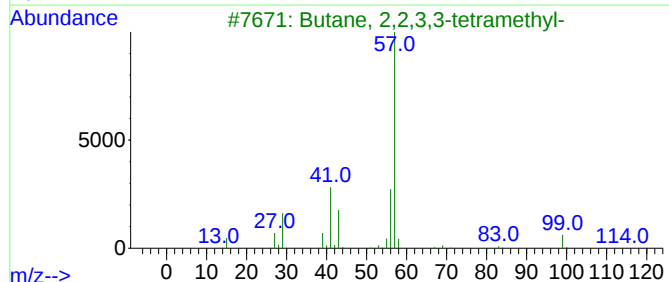
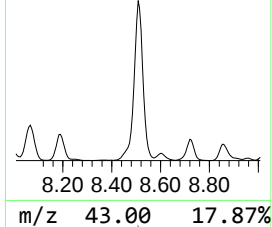
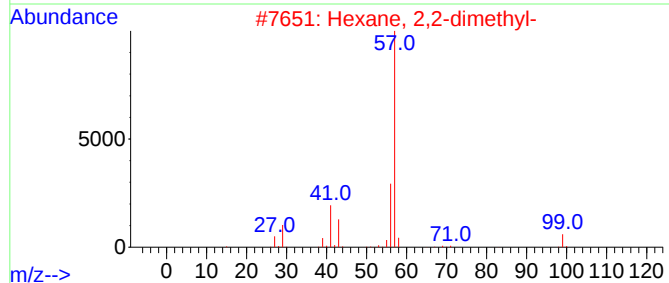
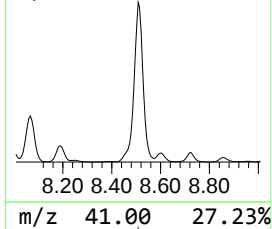
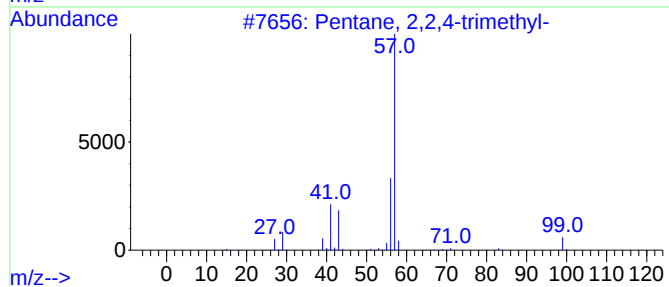
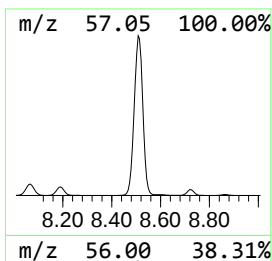
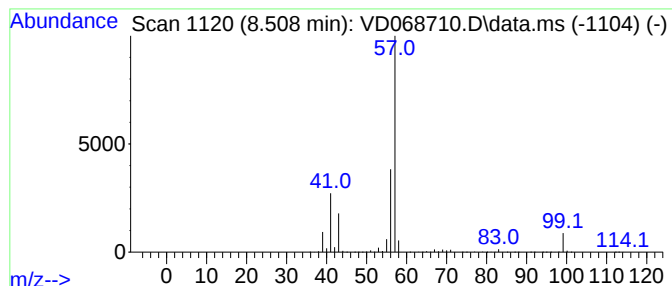
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031921S.M
Quant Title : SW846 8260

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

Peak Number 5 Pentane, 2,2,4-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.508	321.63 ug/l	43375400	1,4-Difluorobenzene	8.867

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	83
2			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	83
3			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
4			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	64
5			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
Data File : VD068710.D
Acq On : 29 Mar 2021 20:50
Operator : VA/SY
Sample : M1836-01
Misc : 8.84G/5.00ml/MSVOA_D/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B1-0-2

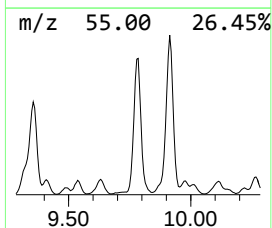
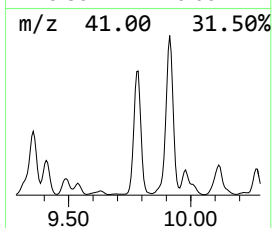
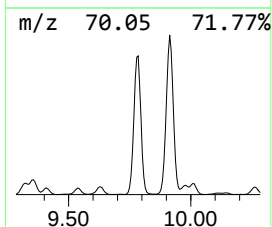
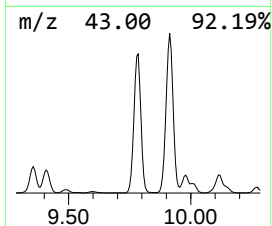
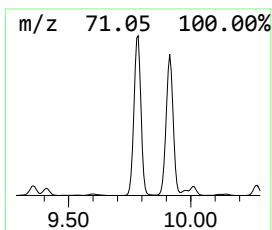
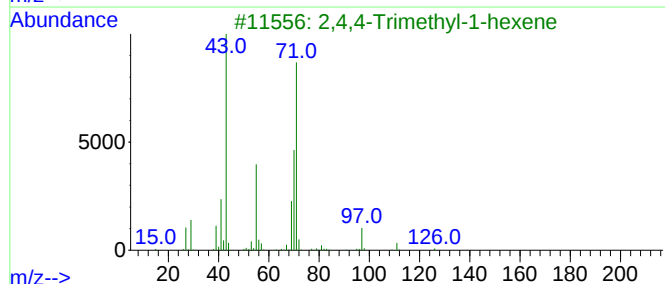
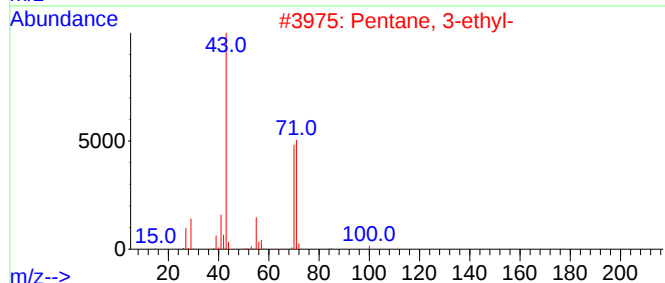
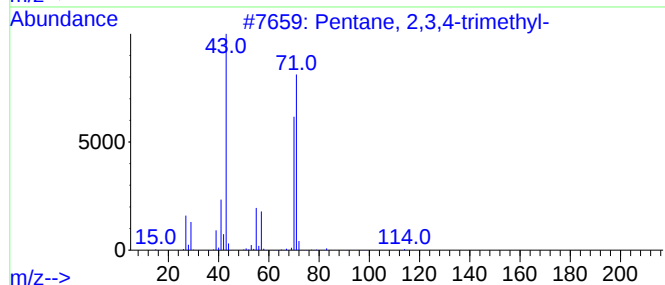
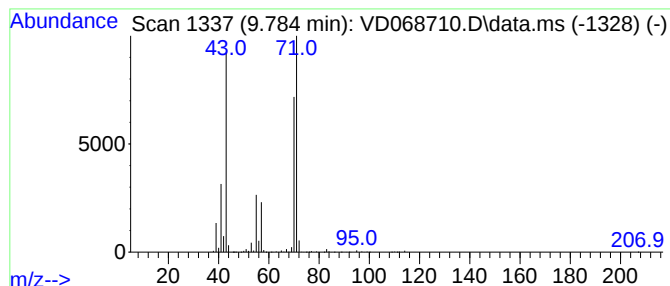
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031921S.M
Quant Title : SW846 8260

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

Peak Number 6 Pentane, 2,3,4-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.784	162.15 ug/l	21867600	1,4-Difluorobenzene	8.867

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2			Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
3			2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	78
4			Hexane, 3-methyl-	100	C7H16	000589-34-4	78
5			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
Data File : VD068710.D
Acq On : 29 Mar 2021 20:50
Operator : VA/SY
Sample : M1836-01
Misc : 8.84G/5.00ml/MSVOA_D/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B1-0-2

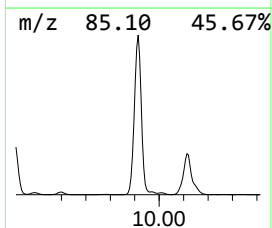
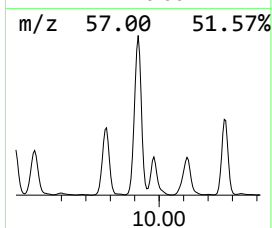
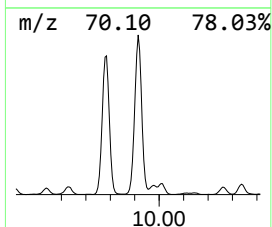
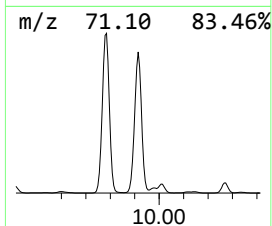
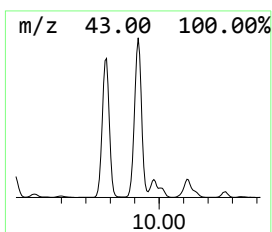
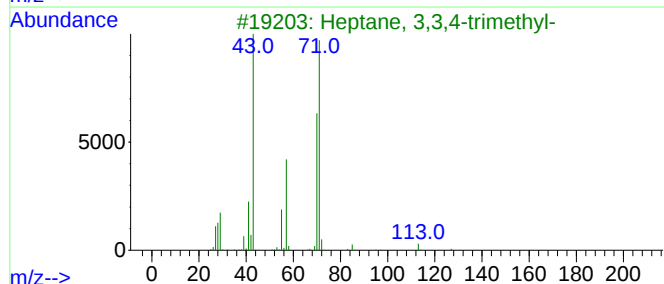
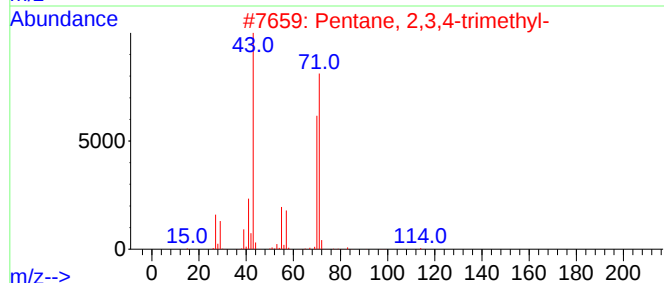
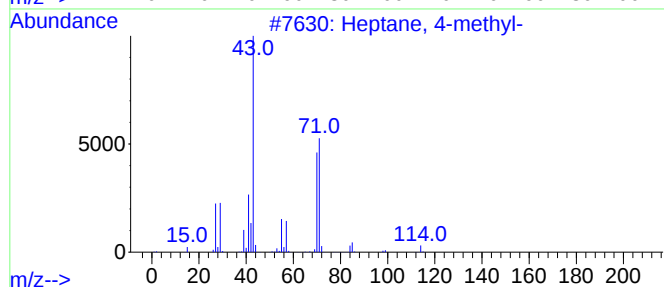
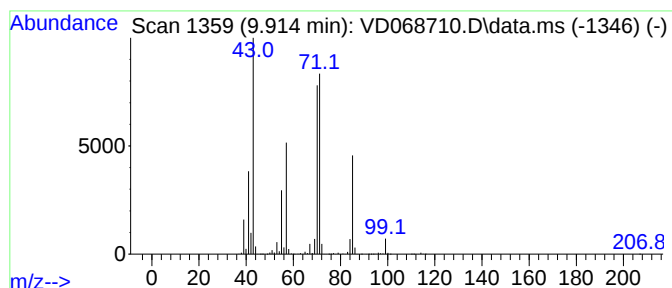
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031921S.M
Quant Title : SW846 8260

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

Peak Number 7 Heptane, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.914	237.79 ug/l	32068500	1,4-Difluorobenzene	8.867

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 4-methyl-	114	C8H18	000589-53-7	83
2			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	72
3			Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	64
4			Pentane, 3-ethyl-	100	C7H16	000617-78-7	59
5			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
Data File : VD068710.D
Acq On : 29 Mar 2021 20:50
Operator : VA/SY
Sample : M1836-01
Misc : 8.84G/5.00ml/MSVOA_D/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B1-0-2

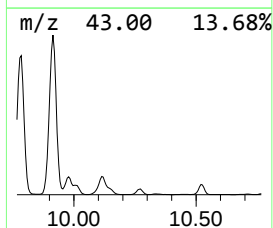
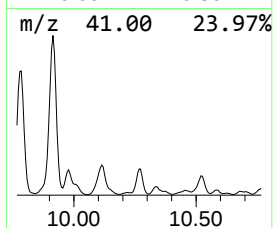
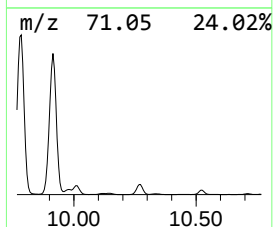
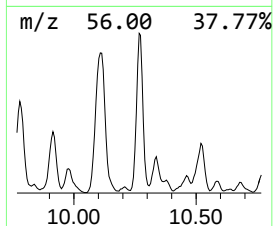
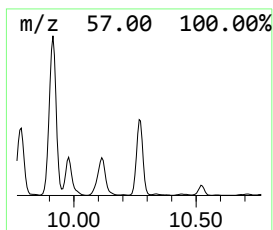
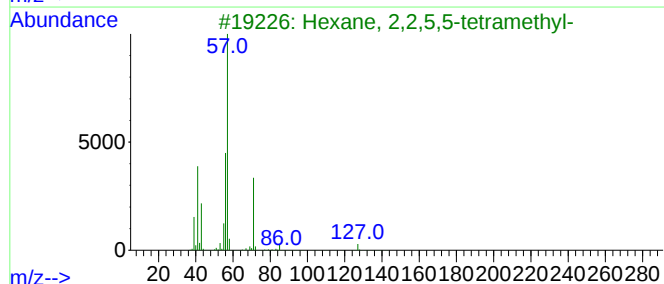
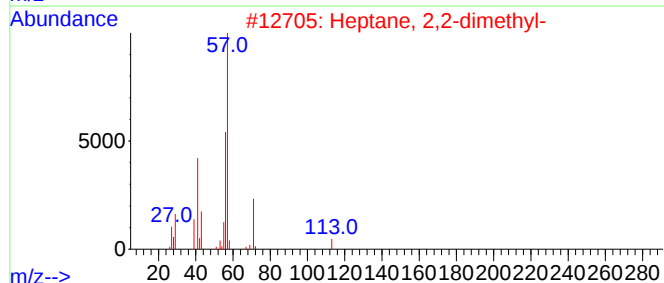
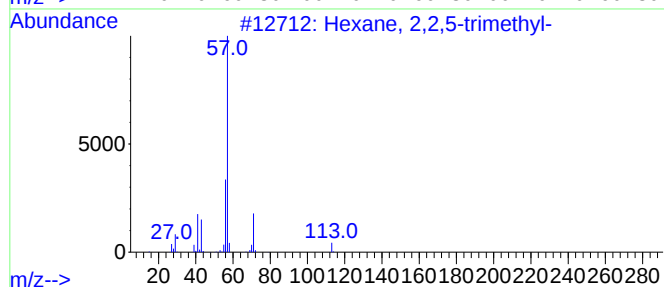
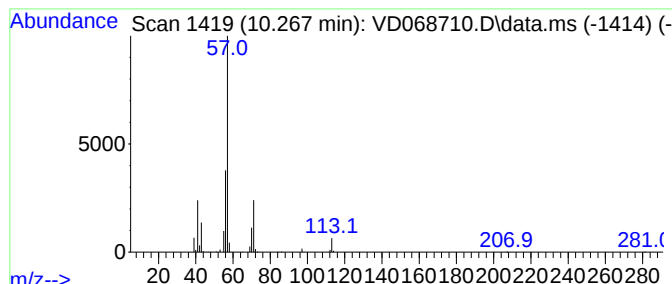
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031921S.M
Quant Title : SW846 8260

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

Peak Number 8 Hexane, 2,2,5-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.267	50.68 ug/l	3336480	Chlorobenzene-d5	11.643

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	72
2			Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	64
3			Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	59
4			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	59
5			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
Data File : VD068710.D
Acq On : 29 Mar 2021 20:50
Operator : VA/SY
Sample : M1836-01
Misc : 8.84G/5.00ml/MSVOA_D/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B1-0-2

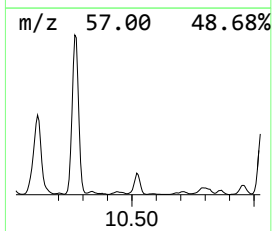
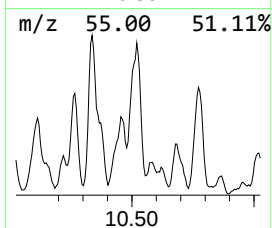
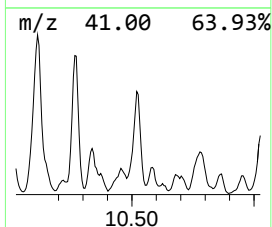
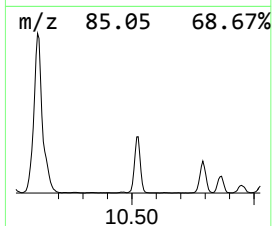
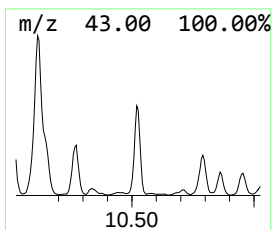
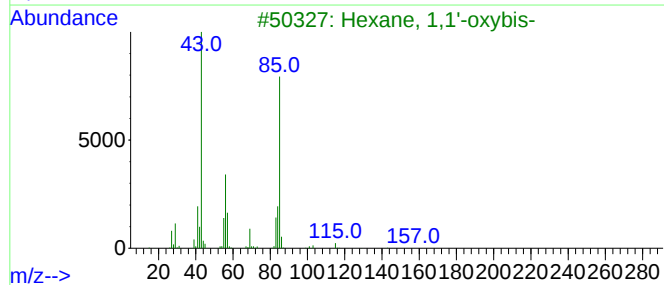
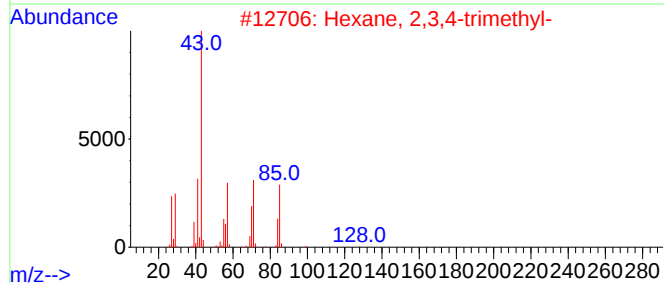
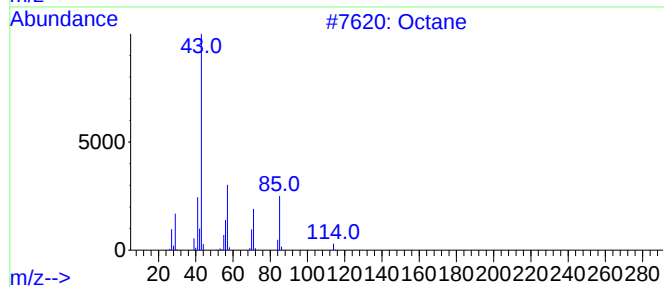
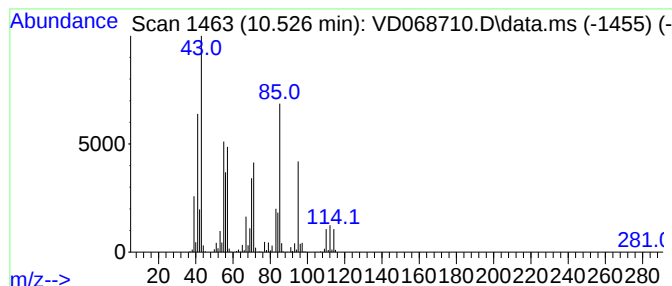
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031921S.M
Quant Title : SW846 8260

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

Peak Number 9 unknown10.526 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.526	45.00 ug/l	2962470	Chlorobenzene-d5	11.643

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane	114	C8H18	000111-65-9	49
2			Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	35
3			Hexane, 1,1'-oxybis-	186	C12H26O	000112-58-3	22
4			1-Decanol, 10-[(tetrahydro-2H-py...	258	C15H30O3	043047-93-4	22
5			1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	22



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
Data File : VD068710.D
Acq On : 29 Mar 2021 20:50
Operator : VA/SY
Sample : M1836-01
Misc : 8.84G/5.00ml/MSVOA_D/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B1-0-2

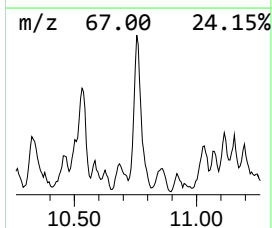
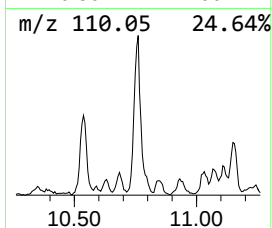
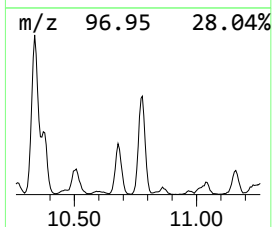
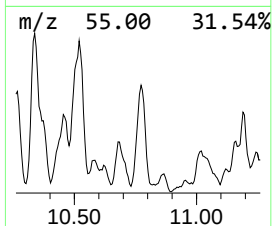
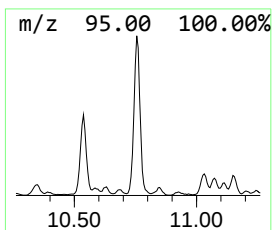
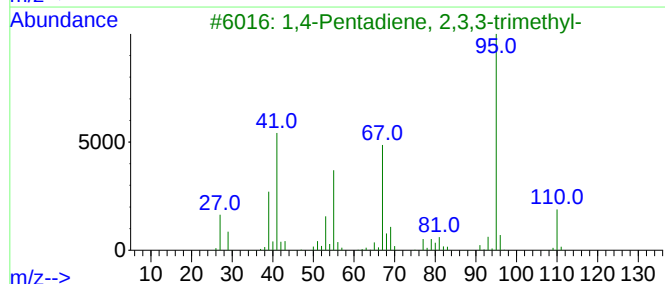
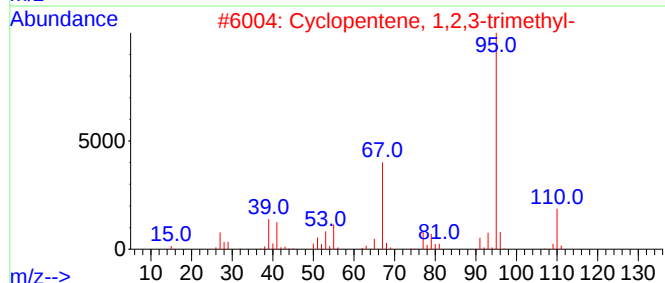
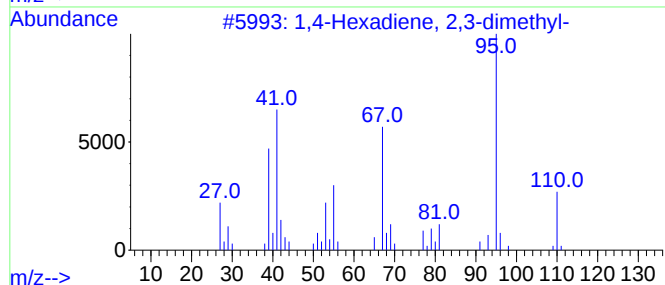
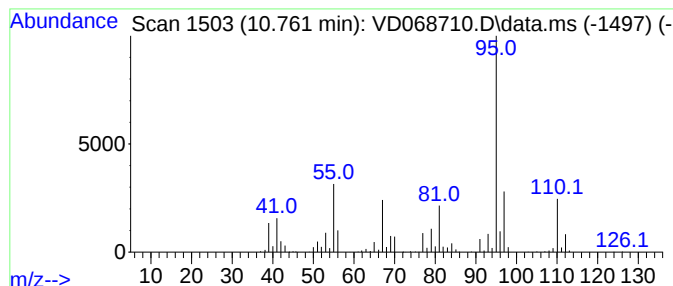
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031921S.M
Quant Title : SW846 8260

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

Peak Number 10 1,4-Hexadiene, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.761	38.47 ug/l	2532500	Chlorobenzene-d5	11.643

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Hexadiene, 2,3-dimethyl-	110	C8H14	018669-52-8	76
2			Cyclopentene, 1,2,3-trimethyl-	110	C8H14	000473-91-6	68
3			1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	59
4			Bicyclo[3.1.0]hexane, 1,5-dimethyl-	110	C8H14	1000142-17-5	59
5			1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD032921\
Data File : VD068710.D
Acq On : 29 Mar 2021 20:50
Operator : VA/SY
Sample : M1836-01
Misc : 8.84G/5.00ml/MSVOA_D/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
B1-0-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031921S.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
Butane, 2-methyl-	3.032	96.6	ug/l	22868000	1	7.979	11834000	50.0
Pentane	3.408	40.2	ug/l	9511710	1	7.979	11834000	50.0
Pentane, 2-methyl-	5.014	57.4	ug/l	13585800	1	7.979	11834000	50.0
Pentane, 3-methyl-	5.491	38.6	ug/l	9146880	1	7.979	11834000	50.0
Pentane, 2,2,4-...	8.508	321.6	ug/l	43375400	2	8.867	6743090	50.0
Pentane, 2,3,4-...	9.784	162.2	ug/l	21867600	2	8.867	6743090	50.0
Heptane, 4-methyl-	9.914	237.8	ug/l	32068500	2	8.867	6743090	50.0
Hexane, 2,2,5-t...	10.267	50.7	ug/l	3336480	3	11.643	3291740	50.0
unknown10.526	10.526	45.0	ug/l	2962470	3	11.643	3291740	50.0
1,4-Hexadiene, ...	10.761	38.5	ug/l	2532500	3	11.643	3291740	50.0