

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD010621\
 Data File : VD068013.D
 Acq On : 06 Jan 2021 15:01
 Operator : VA/SY
 Sample : M1022-02
 Misc : 5.06G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 40967

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\82D122220S.M
 Title : SW846 8260

Signal : TIC: VD068013.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.150	363	379	412	rBV2	21070335	71774435	36.31%	7.608%
2	7.955	1010	1026	1037	rBV	6833709	18443112	9.33%	1.955%
3	8.067	1037	1045	1056	rVV	2559904	6503923	3.29%	0.689%
4	8.191	1056	1066	1083	rVV	10473681	24933620	12.61%	2.643%
5	8.455	1101	1111	1113	rBV2	1813936	3978341	2.01%	0.422%
6	8.508	1113	1120	1130	rVV	7779580	20166056	10.20%	2.138%
7	8.597	1130	1135	1147	rVB	904147	2091842	1.06%	0.222%
8	8.726	1147	1157	1172	rBV	16842527	34302537	17.35%	3.636%
9	9.355	1248	1264	1270	rBV	7919314	18536285	9.38%	1.965%
10	9.408	1270	1273	1280	rVV	2257756	3958050	2.00%	0.420%
11	9.538	1290	1295	1301	rVV	1503727	3025398	1.53%	0.321%
12	9.779	1326	1336	1347	rBV	6257801	12936462	6.54%	1.371%
13	9.914	1347	1359	1365	rBV	9300060	19722484	9.98%	2.091%
14	9.979	1365	1370	1383	rVB2	3111249	7543432	3.82%	0.800%
15	10.120	1383	1394	1407	rVB	2506560	6312272	3.19%	0.669%
16	10.285	1407	1422	1427	rBV	2259760	5767982	2.92%	0.611%
17	10.344	1427	1432	1436	rVV2	3538081	6969739	3.53%	0.739%
18	10.420	1436	1445	1456	rVV4	69637211	197675385	100.00%	20.954%
19	10.526	1456	1463	1470	rVB	7924631	14655264	7.41%	1.554%
20	10.685	1485	1490	1498	rVB	1765951	3269420	1.65%	0.347%
21	10.779	1498	1506	1516	rBV2	2785444	6260370	3.17%	0.664%
22	10.955	1527	1536	1541	rBV	1992553	3425506	1.73%	0.363%
23	11.055	1548	1553	1560	rVB	1427867	2602602	1.32%	0.276%
24	11.126	1560	1565	1568	rBV2	1495073	2418629	1.22%	0.256%
25	11.196	1568	1577	1581	rVV	4068041	8565090	4.33%	0.908%
26	11.243	1581	1585	1592	rVV	3046615	5785319	2.93%	0.613%
27	11.355	1599	1604	1608	rBV2	2147634	3584098	1.81%	0.380%
28	11.432	1608	1617	1629	rVB4	6015676	16528833	8.36%	1.752%
29	11.532	1629	1634	1639	rBV	4940460	7960919	4.03%	0.844%
30	11.749	1663	1671	1680	rBV	20632185	38601626	19.53%	4.092%
31	11.855	1680	1689	1700	rBV3	61237559	137394781	69.51%	14.564%
32	11.938	1700	1703	1708	rVB2	2962475	4553342	2.30%	0.483%
33	12.085	1719	1728	1732	rBV4	1065585	2206461	1.12%	0.234%
34	12.185	1732	1745	1753	rBV	26562343	55903449	28.28%	5.926%
35	12.267	1753	1759	1764	rVB	3460116	5986753	3.03%	0.635%
36	12.355	1769	1774	1776	rBV3	2151246	3339963	1.69%	0.354%

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Peak Location: TOP

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37	12.396	1776	1781	1790	rVB4	3609195	9425142	4.77%	0.999%
38	12.479	1790	1795	1799	rBV2	1823677	3281834	1.66%	0.348%
39	12.667	1824	1827	1832	rVB	1998237	2577434	1.30%	0.273%
40	12.820	1840	1853	1858	rVB2	3487653	9421792	4.77%	0.999%
41	12.885	1858	1864	1867	rBV	9401609	15712970	7.95%	1.666%
42	12.920	1867	1870	1871	rVV	4652033	6037067	3.05%	0.640%
43	12.949	1871	1875	1882	rVV2	15492080	30316093	15.34%	3.214%
44	13.008	1882	1885	1891	rVB2	1879563	3063847	1.55%	0.325%
45	13.126	1896	1905	1911	rBV	3486891	8701838	4.40%	0.922%
46	13.185	1911	1915	1918	rBV2	2682124	4251989	2.15%	0.451%
47	13.267	1924	1929	1935	rBV	10785243	17120774	8.66%	1.815%
48	13.432	1948	1957	1960	rBV2	4002677	8437597	4.27%	0.894%
49	13.614	1983	1988	1999	rVB3	4429529	12958343	6.56%	1.374%
50	13.773	2011	2015	2019	rBV5	3445884	5690534	2.88%	0.603%
51	13.896	2031	2036	2042	rBV2	4964229	7795321	3.94%	0.826%
52	13.973	2045	2049	2055	rVV	1161690	2074032	1.05%	0.220%
53	14.120	2069	2074	2079	rVB	1826806	2855011	1.44%	0.303%
54	14.226	2087	2092	2098	rVB2	1696654	2879952	1.46%	0.305%
55	14.832	2187	2195	2200	rBV3	1383988	3084458	1.56%	0.327%

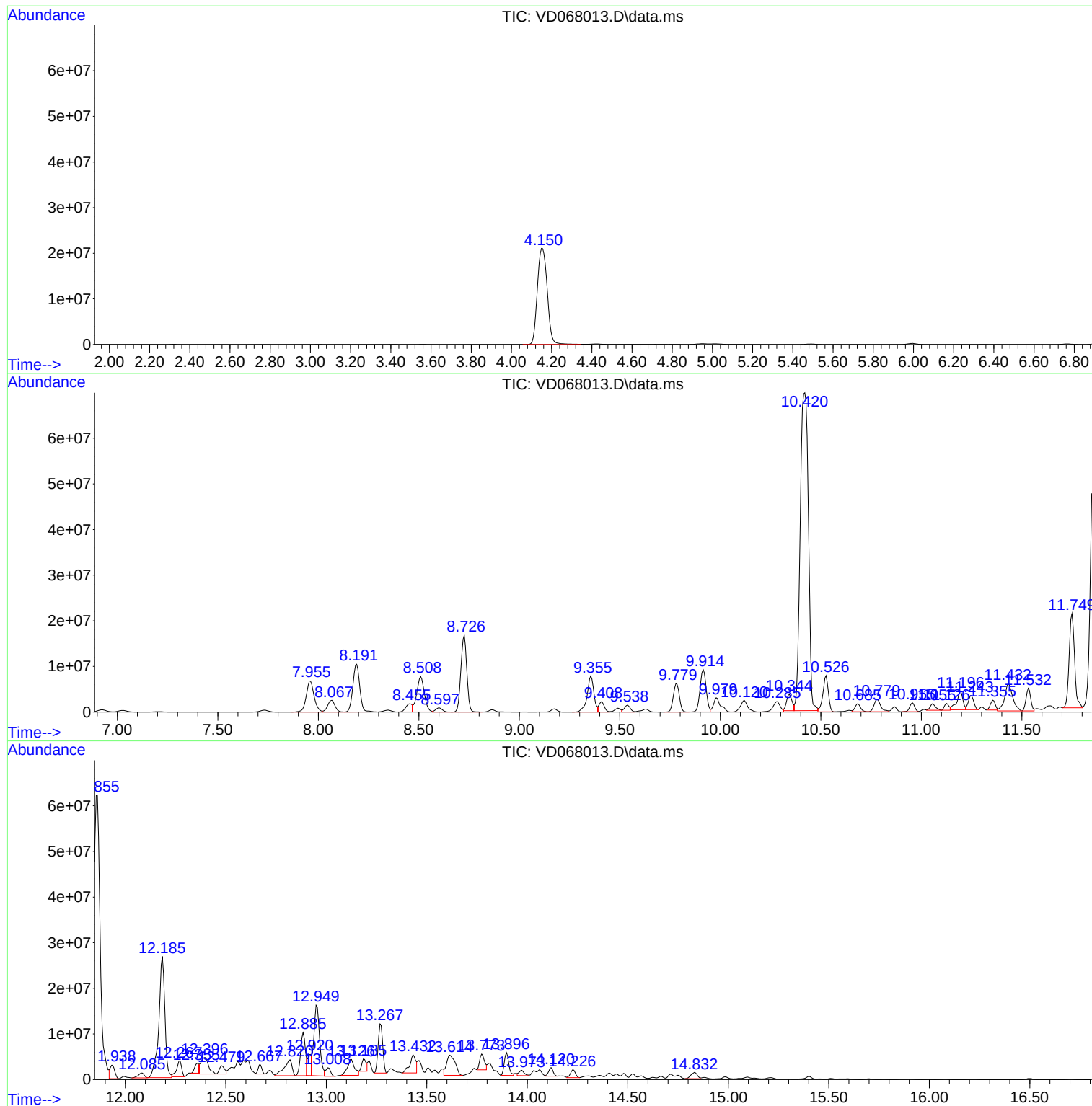
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D122220S.M
Quant Title : SW846 8260

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



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ALS Vial : 6 Sample Multiplier: 1

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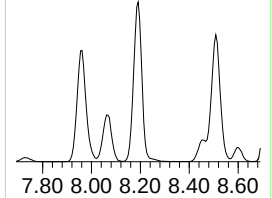
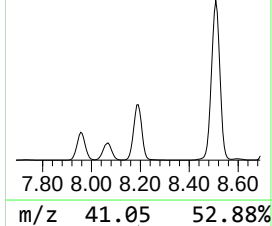
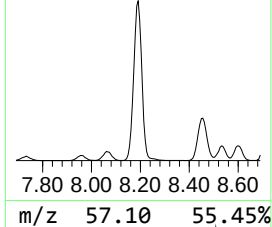
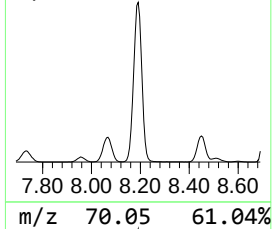
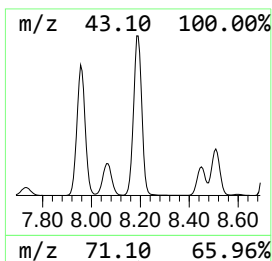
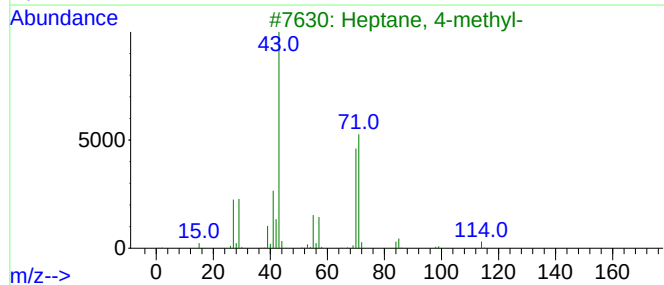
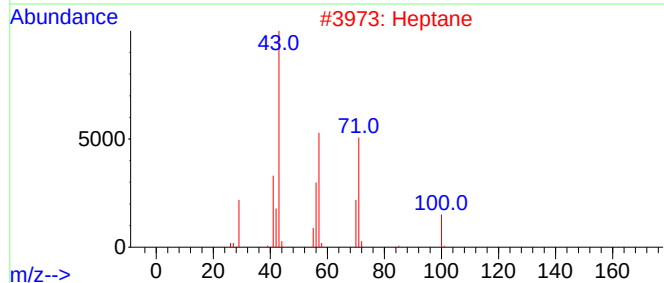
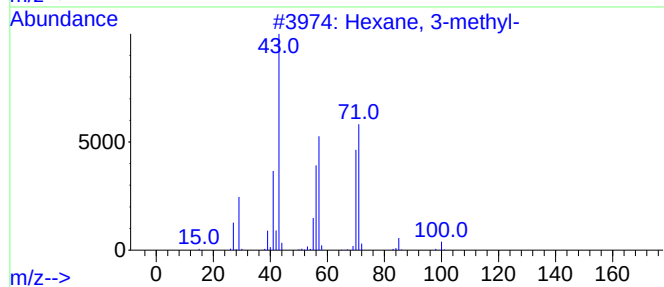
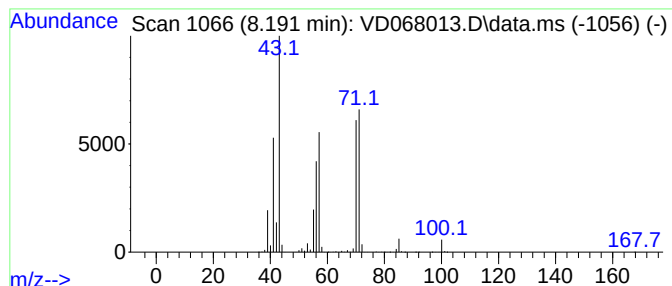
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D12220S.M
Quant Title : SW846 8260

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

Peak Number 1 Hexane, 3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.191	67.60 ug/l	24933600	Pentafluorobenzene	7.979

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	94
2			Heptane	100	C7H16	000142-82-5	80
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53
5			Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	53



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Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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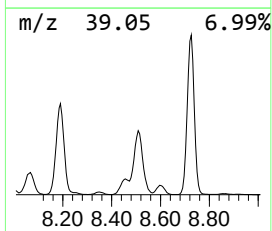
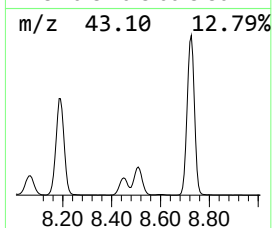
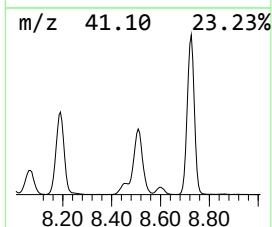
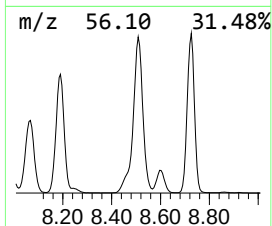
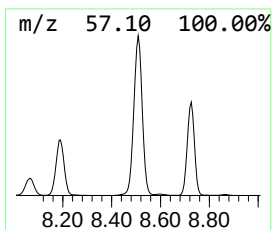
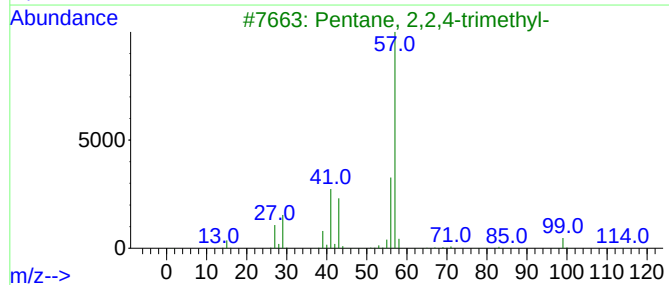
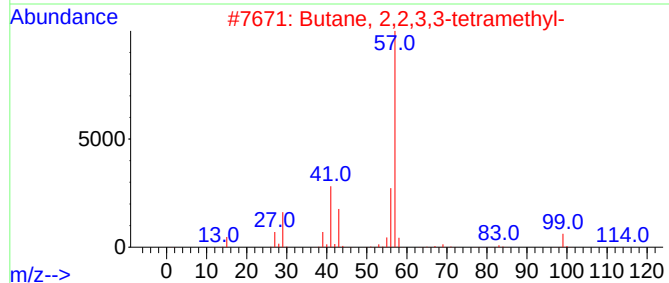
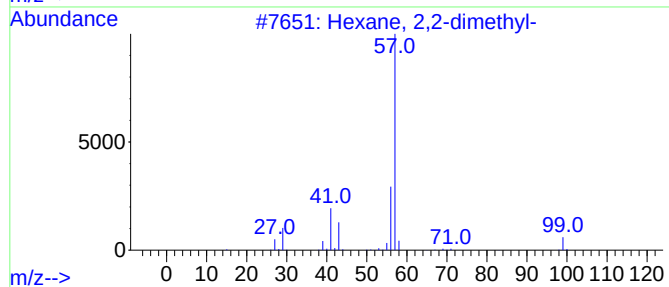
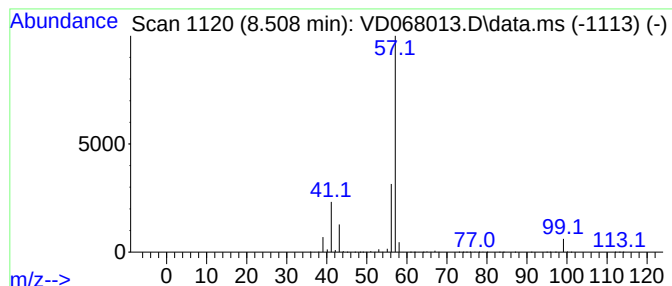
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D122220S.M
Quant Title : SW846 8260

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

Peak Number 2 Hexane, 2,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.508	29.39 ug/l	20166100	1,4-Difluorobenzene	8.867

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	78
2			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	64
3			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	64
4			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	39
5			Acetamide, N-2-propenyl-	99	C5H9NO	000692-33-1	28



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD010621\
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Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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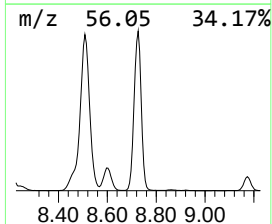
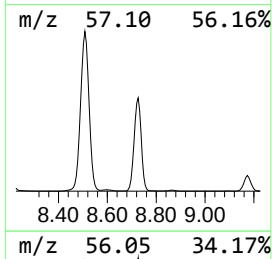
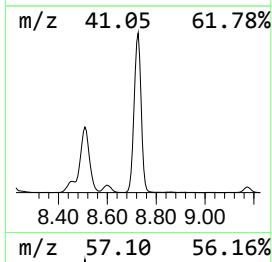
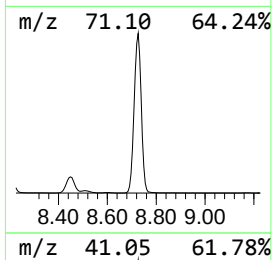
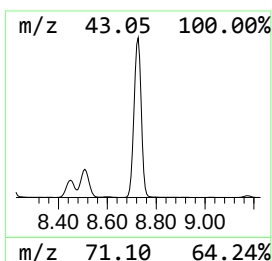
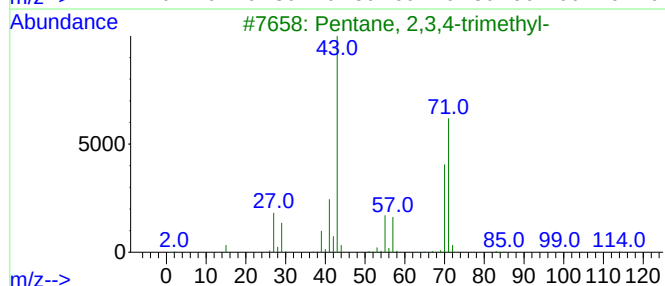
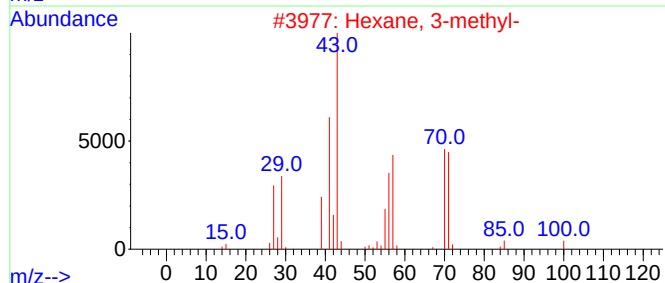
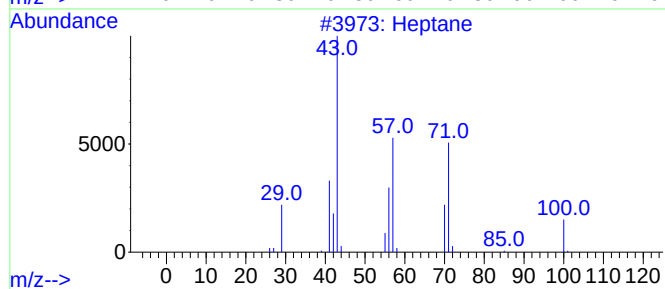
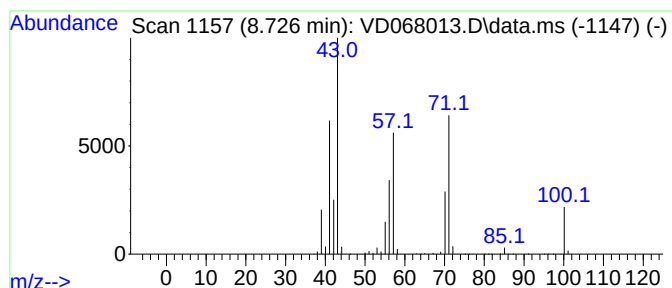
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D12220S.M
Quant Title : SW846 8260

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

Peak Number 3 Heptane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.726	50.00 ug/l	34302500	1,4-Difluorobenzene	8.867

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	40
3			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	40
4			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38
5			Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	37



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Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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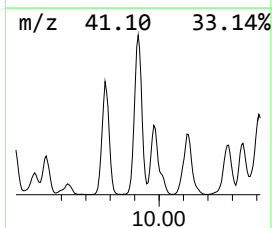
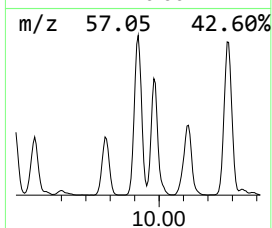
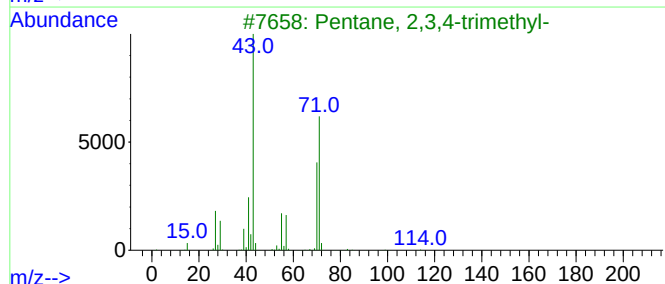
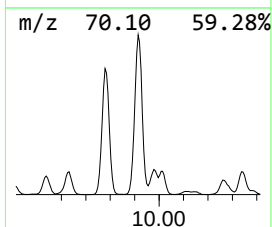
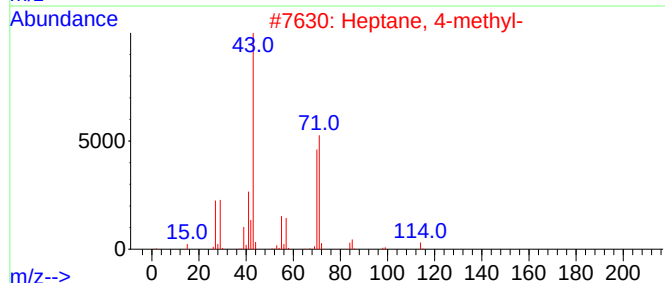
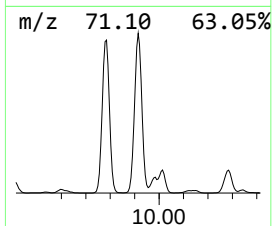
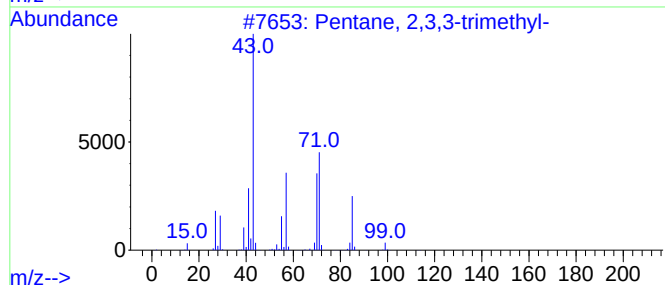
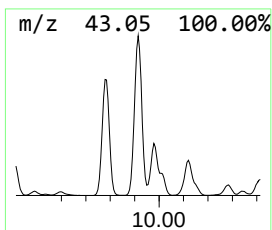
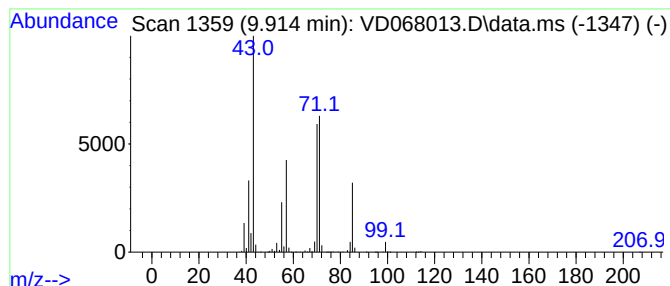
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D12220S.M
Quant Title : SW846 8260

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

Peak Number 4 Pentane, 2,3,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.914	28.75 ug/l	19722500	1,4-Difluorobenzene	8.867

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	80
2			Heptane, 4-methyl-	114	C8H18	000589-53-7	64
3			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	64
4			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	59
5			Hexane, 3-methyl-	100	C7H16	000589-34-4	59



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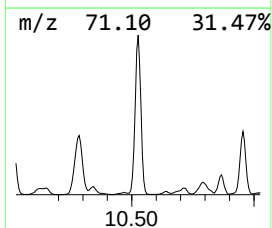
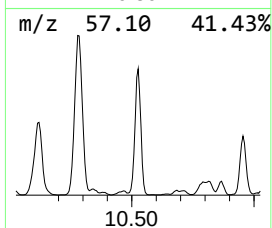
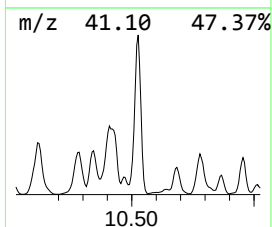
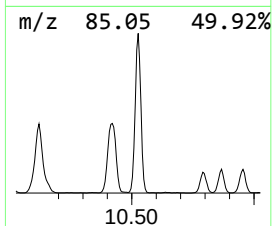
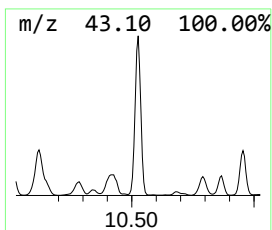
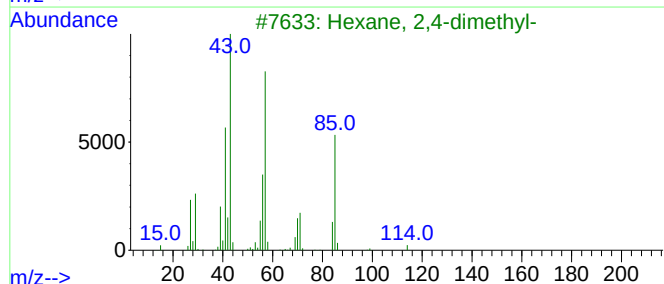
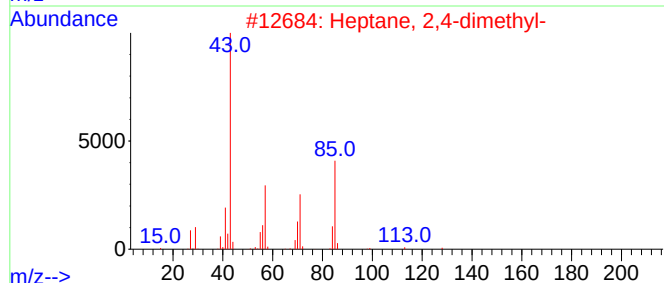
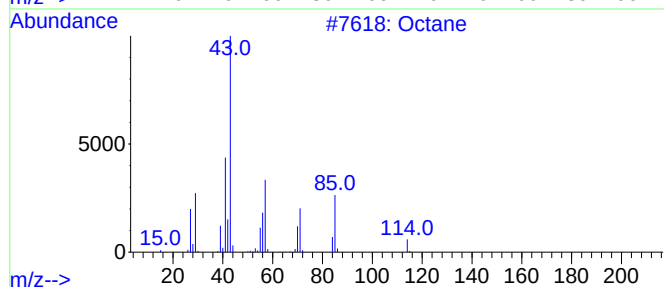
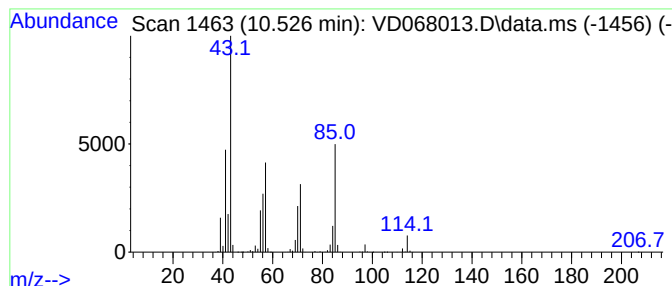
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Quant Title : SW846 8260

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

Peak Number 5 Octane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.526	18.98 ug/l	14655300	Chlorobenzene-d5	11.649

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane	114	C8H18	000111-65-9	91
2			Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	80
3			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	64
4			Decane, 5,6-dimethyl-	170	C12H26	001636-43-7	59
5			Oxalic acid, butyl isohexyl ester	230	C12H22O4	1000309-32-7	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD010621\
Data File : VD068013.D
Acq On : 06 Jan 2021 15:01
Operator : VA/SY
Sample : M1022-02
Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
40967

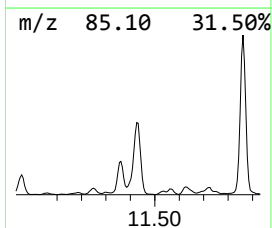
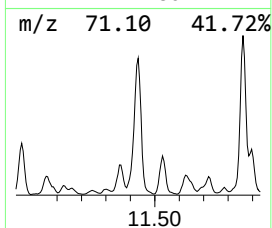
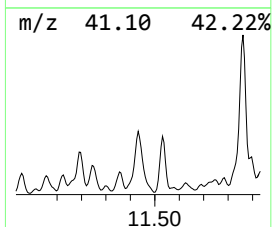
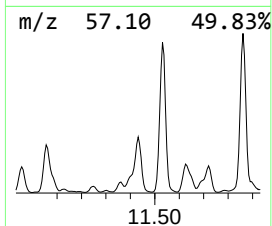
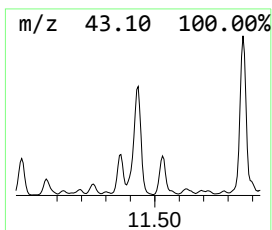
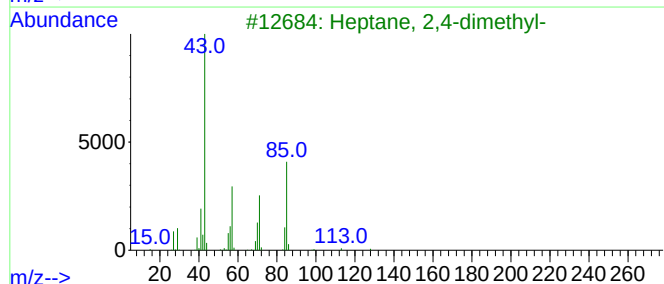
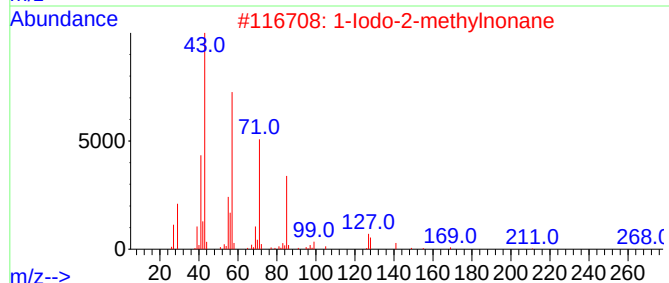
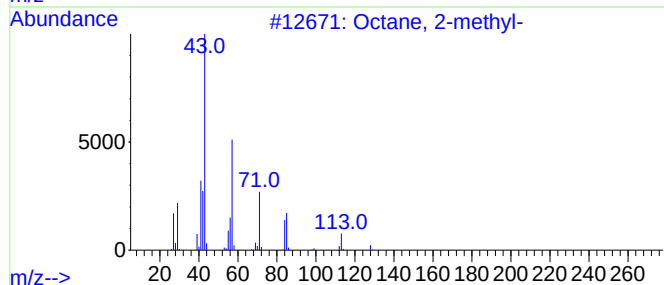
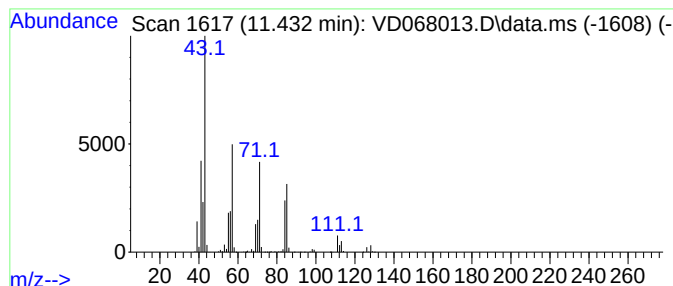
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Quant Title : SW846 8260

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

Peak Number 6 Octane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.432	21.41 ug/l	16528800	Chlorobenzene-d5	11.649

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2-methyl-	128	C9H20	003221-61-2	74
2			1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	53
3			Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	53
4			Decane, 2,4,6-trimethyl-	184	C13H28	062108-27-4	50
5			Hexane, 2,3,5-trimethyl-	128	C9H20	001069-53-0	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD010621\
Data File : VD068013.D
Acq On : 06 Jan 2021 15:01
Operator : VA/SY
Sample : M1022-02
Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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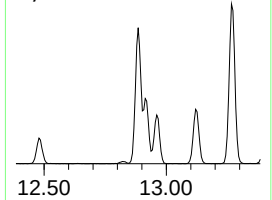
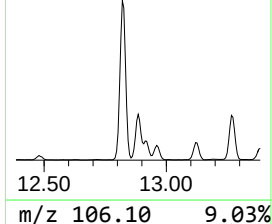
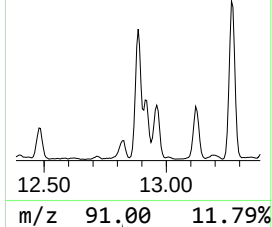
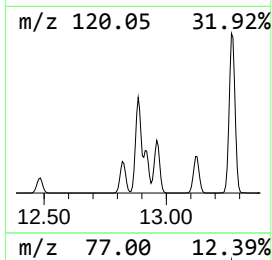
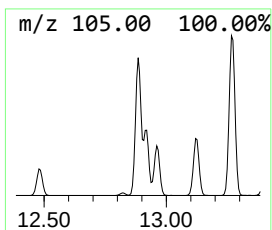
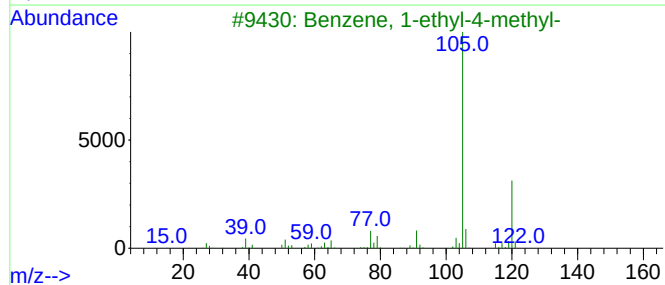
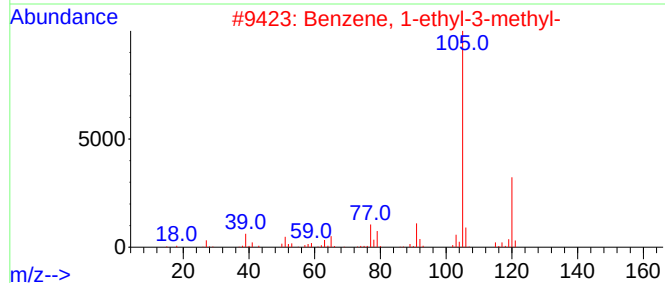
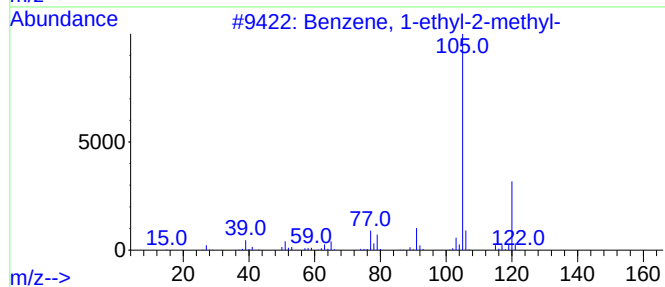
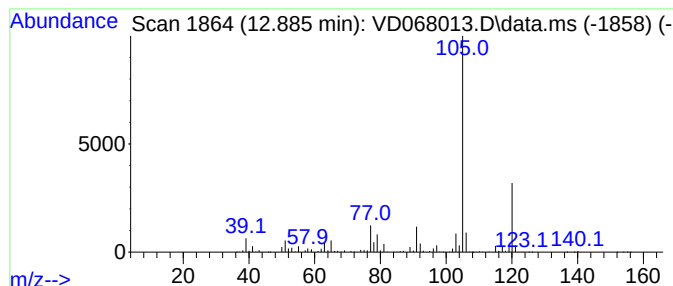
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.885	60.63 ug/l	15713000	1,4-Dichlorobenzene-d4	13.579

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	95
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD010621\
Data File : VD068013.D
Acq On : 06 Jan 2021 15:01
Operator : VA/SY
Sample : M1022-02
Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
40967

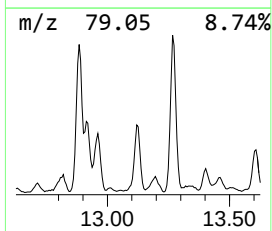
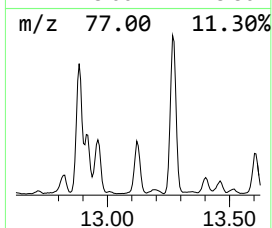
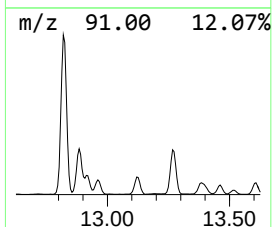
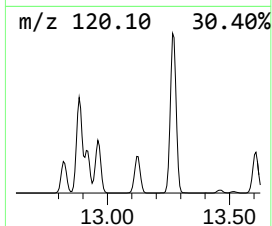
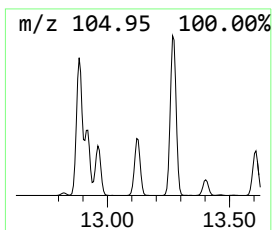
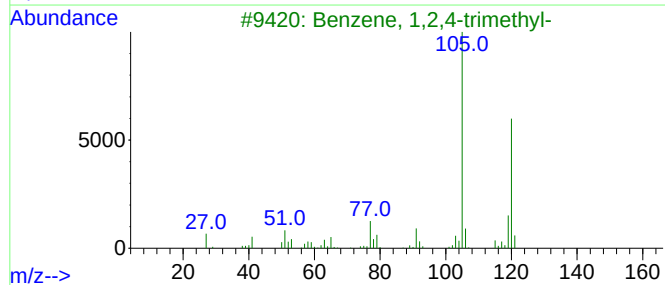
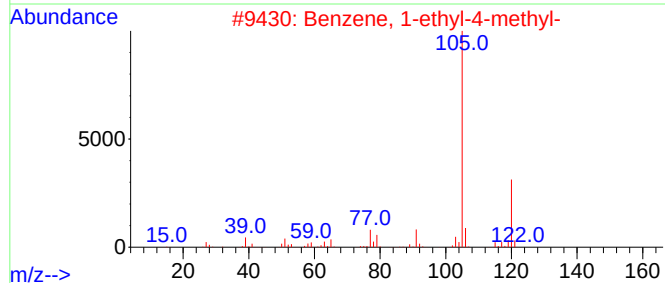
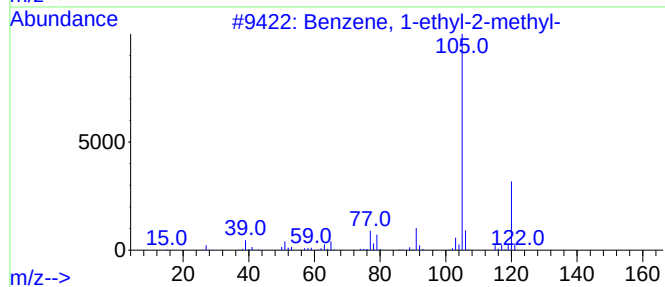
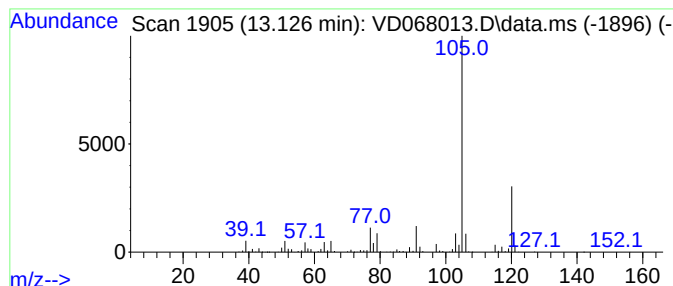
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.126	33.58 ug/l	8701840	1,4-Dichlorobenzene-d4	13.579

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5			Mesitylene	120	C9H12	000108-67-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD010621\
Data File : VD068013.D
Acq On : 06 Jan 2021 15:01
Operator : VA/SY
Sample : M1022-02
Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
40967

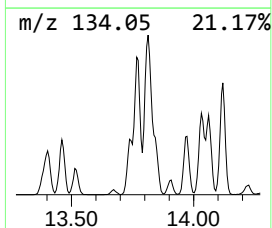
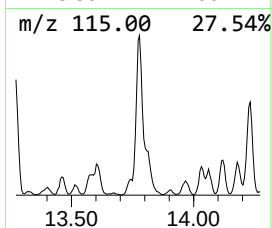
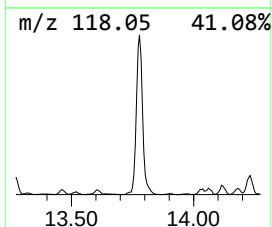
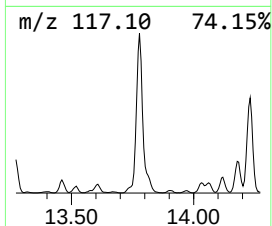
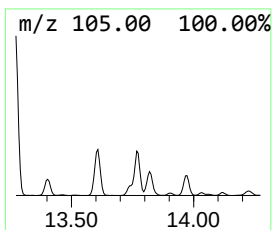
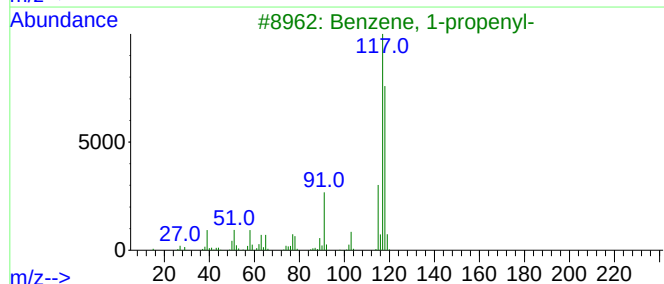
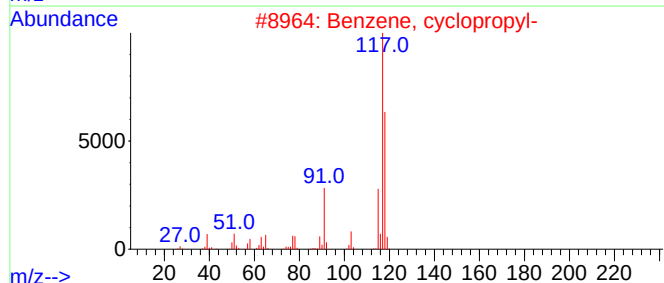
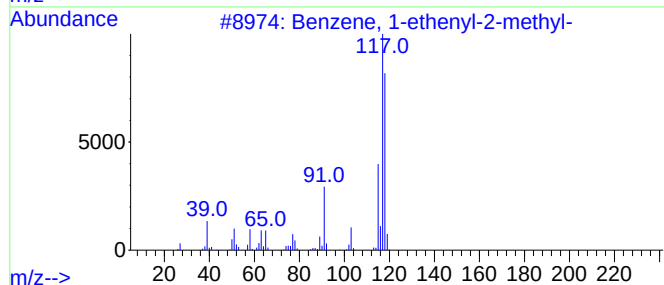
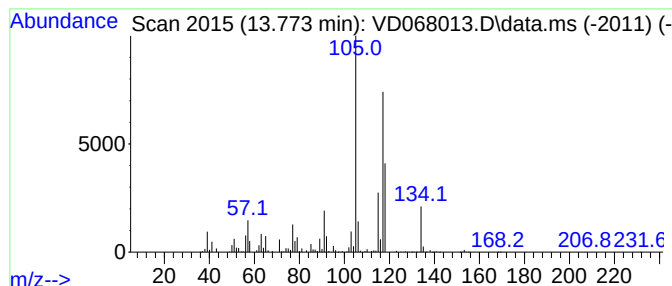
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethenyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.773	21.96 ug/l	5690530	1,4-Dichlorobenzene-d4	13.579

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	92
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
3			Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
4			Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
5			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD010621\
Data File : VD068013.D
Acq On : 06 Jan 2021 15:01
Operator : VA/SY
Sample : M1022-02
Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
40967

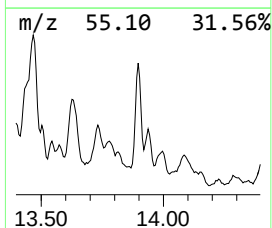
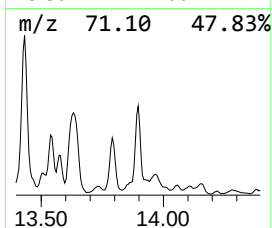
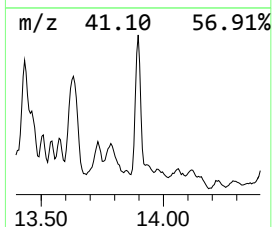
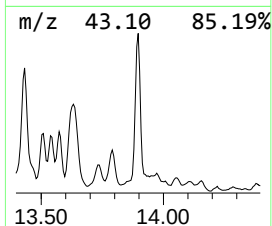
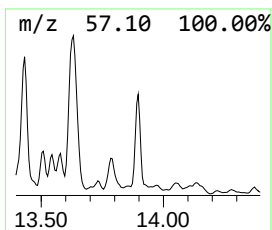
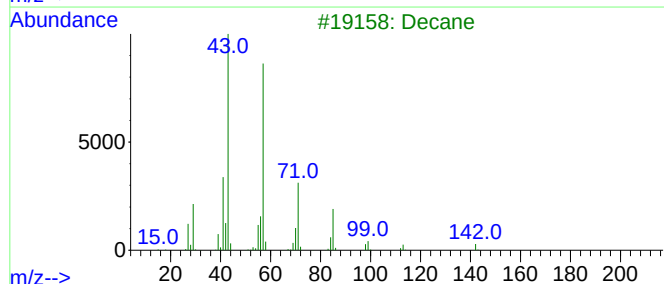
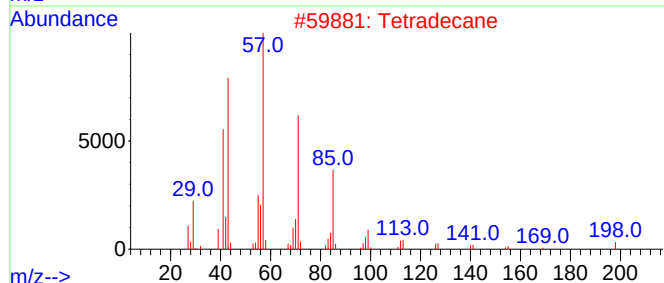
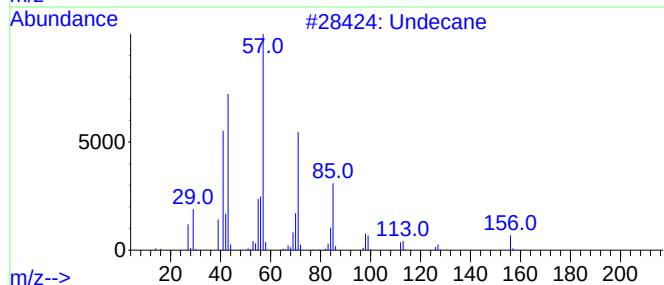
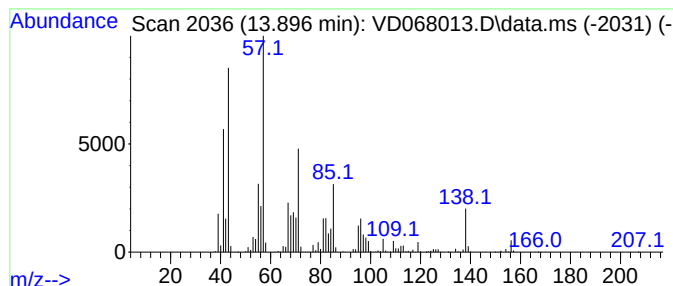
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Quant Title : SW846 8260

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

Peak Number 10 Undecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.896	30.08 ug/l	7795320	1,4-Dichlorobenzene-d4	13.579

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	83
2	Tetradecane			198	C14H30	000629-59-4	47
3	Decane			142	C10H22	000124-18-5	47
4	Tridecane			184	C13H28	000629-50-5	43
5	Hexadecane			226	C16H34	000544-76-3	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD010621\
 Data File : VD068013.D
 Acq On : 06 Jan 2021 15:01
 Operator : VA/SY
 Sample : M1022-02
 Misc : 5.06G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D122220S.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
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Hexane, 2,2-dim...	8.508	29.4	ug/l	20166100	2	8.867	34302500	50.0
Heptane	8.726	50.0	ug/l	34302500	2	8.867	34302500	50.0
Pentane, 2,3,3-...	9.914	28.8	ug/l	19722500	2	8.867	34302500	50.0
Octane	10.526	19.0	ug/l	14655300	3	11.649	38601600	50.0
Octane, 2-methyl-	11.432	21.4	ug/l	16528800	3	11.649	38601600	50.0
Benzene, 1-ethy...	12.885	60.6	ug/l	15713000	4	13.579	12958300	50.0
Benzene, 1-ethy...	13.126	33.6	ug/l	8701840	4	13.579	12958300	50.0
Benzene, 1-ethe...	13.773	22.0	ug/l	5690530	4	13.579	12958300	50.0
Undecane	13.896	30.1	ug/l	7795320	4	13.579	12958300	50.0