

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070722\
 Data File : VD073817.D
 Acq On : 07 Jul 2022 19:28
 Operator : VA/SY
 Sample : N3593-10
 Misc : 4.77G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 20220486A-4A

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\82D070622S.M
 Title : SW846 8260

Signal : TIC: VD073817.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.950	3	5	9	rVB	19571	20190	2.54%	0.226%
2	2.309	61	66	69	rBV2	17928	28305	3.57%	0.318%
3	3.391	244	250	260	rVB6	3793	11070	1.39%	0.124%
4	4.156	370	380	389	rBV	26106	75096	9.46%	0.842%
5	4.226	389	392	399	rVB2	5106	10173	1.28%	0.114%
6	4.397	410	421	432	rBV2	29415	82942	10.45%	0.930%
7	4.950	504	515	524	rBV4	11253	38844	4.89%	0.436%
8	7.203	888	898	906	rBV2	10114	26940	3.39%	0.302%
9	7.908	1008	1018	1023	rBV2	101547	250534	31.56%	2.811%
10	7.967	1023	1028	1038	rVV2	149300	336413	42.38%	3.774%
11	8.061	1038	1044	1055	rVB5	11611	26500	3.34%	0.297%
12	8.185	1058	1065	1070	rVB7	6065	13808	1.74%	0.155%
13	8.320	1080	1088	1099	rBV2	100311	228651	28.81%	2.565%
14	8.497	1111	1118	1129	rVB	37703	94408	11.89%	1.059%
15	8.855	1169	1179	1191	rBV	260365	532909	67.14%	5.978%
16	9.085	1212	1218	1224	rBV3	4893	8102	1.02%	0.091%
17	9.349	1251	1263	1268	rBV4	27445	72709	9.16%	0.816%
18	9.402	1268	1272	1278	rVB4	11180	20954	2.64%	0.235%
19	9.620	1299	1309	1314	rVB6	8570	20907	2.63%	0.235%
20	9.767	1327	1334	1343	rVB	40051	81351	10.25%	0.913%
21	9.908	1346	1358	1364	rBV2	46150	103581	13.05%	1.162%
22	10.091	1383	1389	1394	rBV4	8254	18552	2.34%	0.208%
23	10.255	1405	1417	1423	rBV2	27473	60157	7.58%	0.675%
24	10.332	1423	1430	1443	rVB	431220	793736	100.00%	8.904%
25	10.508	1453	1460	1467	rVB10	9604	27734	3.49%	0.311%
26	10.632	1471	1481	1484	rVV3	9001	18769	2.36%	0.211%
27	10.667	1484	1487	1493	rVV3	8922	17874	2.25%	0.201%
28	10.773	1496	1505	1515	rVV9	15634	47866	6.03%	0.537%
29	10.849	1515	1518	1524	rVB3	7588	12932	1.63%	0.145%
30	10.943	1529	1534	1541	rVB2	20551	39574	4.99%	0.444%
31	11.043	1543	1551	1559	rVV4	16235	42185	5.31%	0.473%
32	11.114	1559	1563	1567	rVV3	10488	16361	2.06%	0.184%
33	11.179	1571	1574	1578	rVV4	12319	22810	2.87%	0.256%
34	11.238	1578	1584	1589	rVV	37714	74747	9.42%	0.839%
35	11.291	1589	1593	1596	rVV5	12761	19397	2.44%	0.218%
36	11.338	1596	1601	1606	rVV5	16673	30961	3.90%	0.347%

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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37	11.385	1606	1609	1614	rVV4	8229	15869	2.00%	0.178%
38	11.443	1614	1619	1627	rVB3	19801	41401	5.22%	0.464%
39	11.520	1627	1632	1635	rBV4	6448	11736	1.48%	0.132%
40	11.632	1644	1651	1661	rBV	370184	671556	84.61%	7.534%
41	11.773	1670	1675	1678	rBV4	9982	16283	2.05%	0.183%
42	11.826	1678	1684	1688	rBV4	21790	43886	5.53%	0.492%
43	11.879	1689	1693	1698	rVV5	23728	40563	5.11%	0.455%
44	11.985	1705	1711	1712	rBV3	18393	29270	3.69%	0.328%
45	12.073	1722	1726	1730	rVB3	14884	21875	2.76%	0.245%
46	12.138	1730	1737	1749	rBV9	51690	172612	21.75%	1.936%
47	12.255	1749	1757	1762	rVB2	73229	154164	19.42%	1.729%
48	12.367	1762	1776	1783	rBV4	117723	361121	45.50%	4.051%
49	12.461	1783	1792	1797	rBV8	44149	117622	14.82%	1.320%
50	12.502	1797	1799	1808	rVB4	34680	69317	8.73%	0.778%
51	12.620	1811	1819	1825	rVB3	359248	647594	81.59%	7.265%
52	12.708	1825	1834	1839	rBV6	66199	168338	21.21%	1.888%
53	12.785	1843	1847	1853	rVB6	79736	151627	19.10%	1.701%
54	12.867	1853	1861	1865	rBV5	60971	154545	19.47%	1.734%
55	12.949	1867	1875	1887	rVB10	79144	304043	38.31%	3.411%
56	13.055	1888	1893	1896	rBV2	65658	125306	15.79%	1.406%
57	13.132	1902	1906	1909	rVB4	29642	42750	5.39%	0.480%
58	13.173	1909	1913	1915	rBV5	40309	62643	7.89%	0.703%
59	13.261	1924	1928	1931	rVB4	24767	36327	4.58%	0.408%
60	13.296	1931	1934	1939	rBV3	44479	67089	8.45%	0.753%
61	13.349	1940	1943	1946	rVV2	24058	24683	3.11%	0.277%
62	13.420	1951	1955	1959	rBV3	56290	99252	12.50%	1.113%
63	13.485	1964	1966	1971	rVB4	30307	35147	4.43%	0.394%
64	13.561	1972	1979	1984	rBV	333907	580499	73.14%	6.512%
65	13.702	1997	2003	2008	rBV6	41062	83252	10.49%	0.934%
66	13.767	2009	2014	2019	rVB8	19130	46736	5.89%	0.524%
67	13.885	2029	2034	2038	rBV5	62690	105750	13.32%	1.186%
68	13.985	2048	2051	2054	rBV5	21499	32603	4.11%	0.366%
69	14.208	2084	2089	2093	rBV6	33737	62493	7.87%	0.701%
70	14.273	2095	2100	2106	rVB6	44074	83509	10.52%	0.937%
71	14.326	2106	2109	2113	rBV5	17312	29522	3.72%	0.331%
72	14.390	2116	2120	2124	rBV4	57587	88108	11.10%	0.988%
73	14.502	2137	2139	2143	rBV5	15164	25136	3.17%	0.282%
74	14.549	2144	2147	2152	rVB4	42550	65617	8.27%	0.736%
75	14.673	2164	2168	2172	rBV5	15080	27771	3.50%	0.312%
76	14.808	2185	2191	2195	rBV7	50893	104707	13.19%	1.175%

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ClientSampleId :
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Integration Parameters: RTEINT.P

Integrator: RTE

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77	14.855	2195	2199	2205	rVB4	60966	114756	14.46%	1.287%
78	15.067	2232	2235	2241	rVB7	32369	49638	6.25%	0.557%
79	15.343	2277	2282	2289	rVB4	55069	106151	13.37%	1.191%
80	15.414	2290	2294	2301	rBV9	26523	67735	8.53%	0.760%
81	15.590	2320	2324	2331	rBV8	23842	47929	6.04%	0.538%
82	15.755	2348	2352	2360	rVB6	53805	102596	12.93%	1.151%
83	16.320	2445	2448	2457	rVB4	33378	74792	9.42%	0.839%

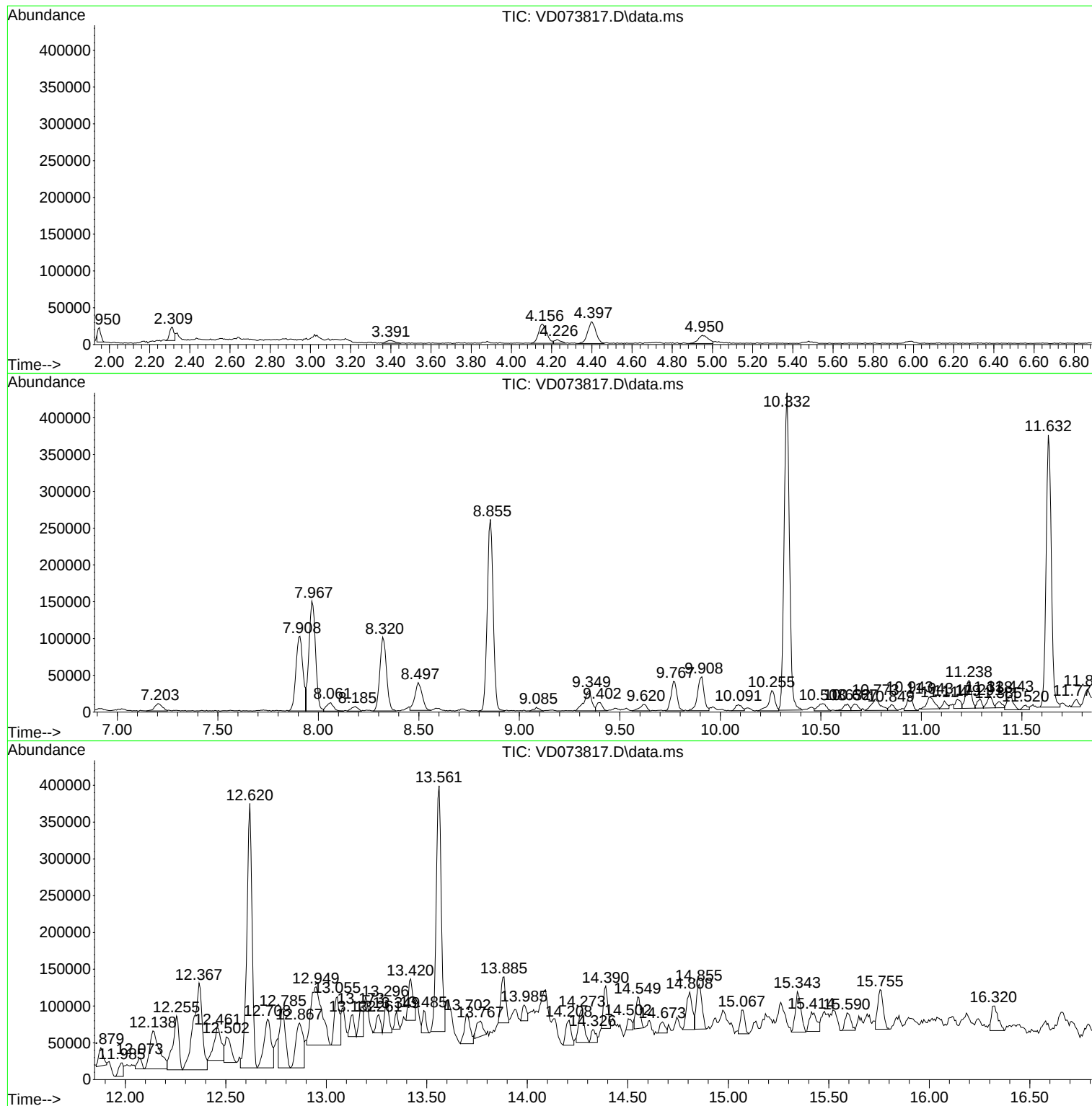
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D070622S.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_D
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20220486A-4A

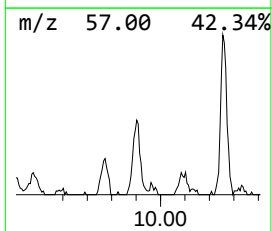
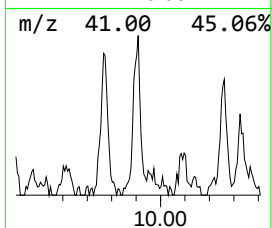
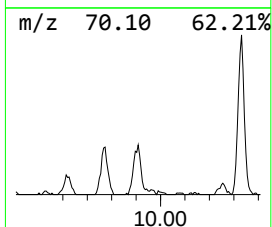
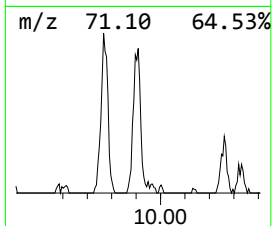
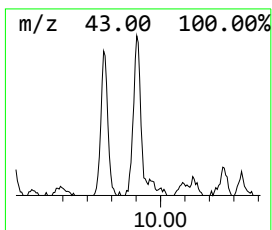
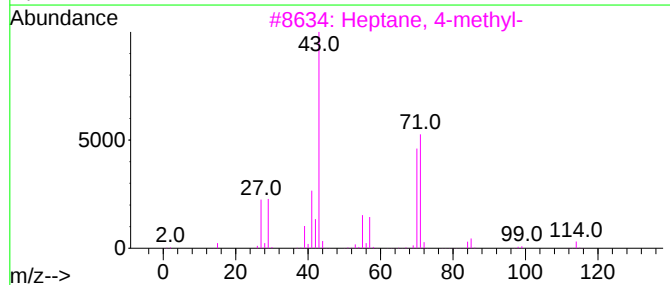
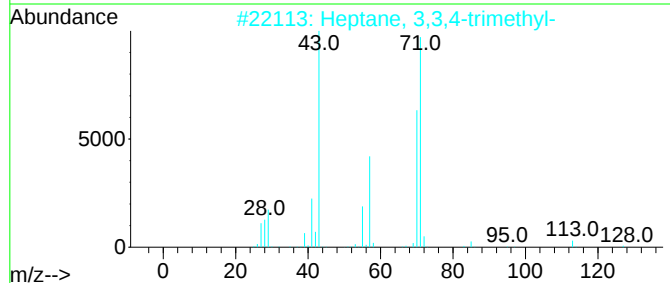
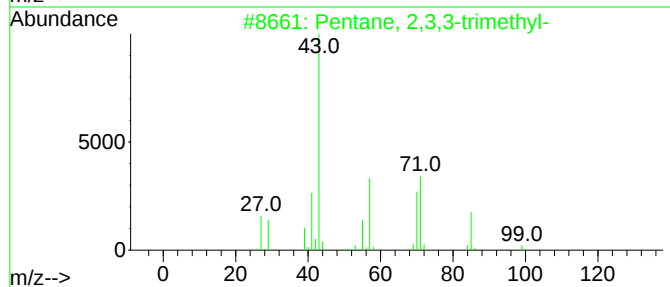
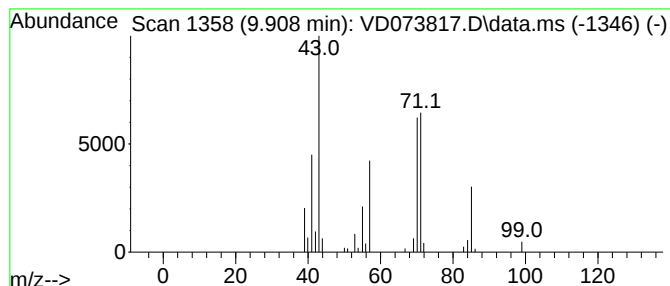
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D070622S.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2,3,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.908	9.72 ug/l	103581	1,4-Difluorobenzene	8.855

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	74
2			Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	64
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	64
4			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64
5			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	59



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Misc : 4.77G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

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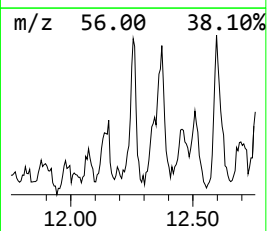
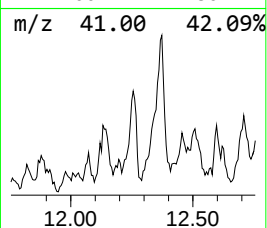
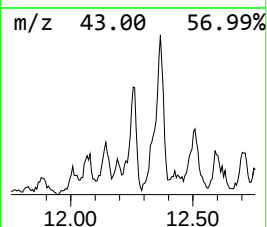
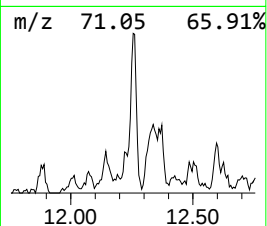
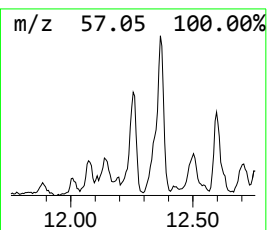
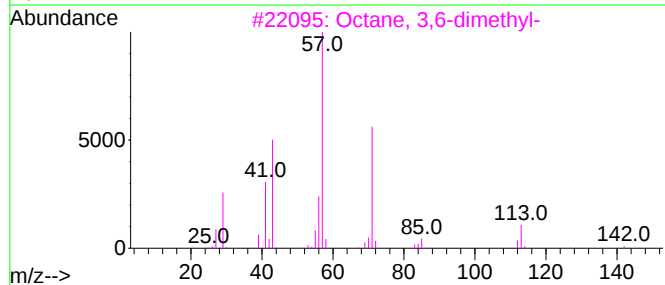
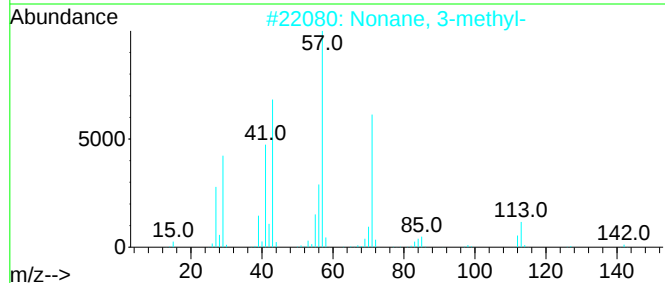
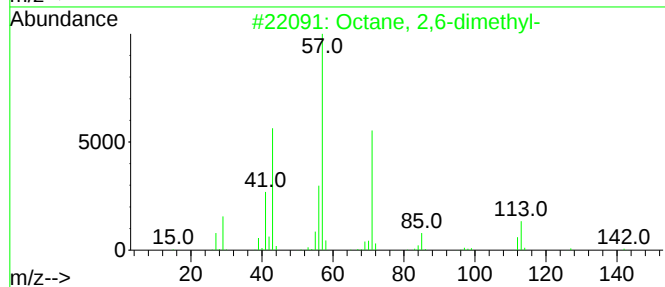
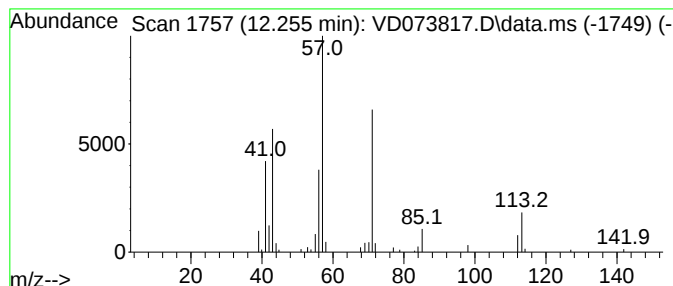
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D070622S.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.255	11.48 ug/l	154164	Chlorobenzene-d5	11.632

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	93
2			Nonane, 3-methyl-	142	C10H22	005911-04-6	90
3			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	64
4			Decane, 3-methyl-	156	C11H24	013151-34-3	53
5			Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	50



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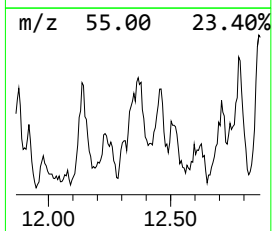
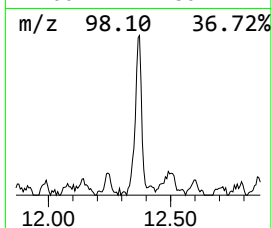
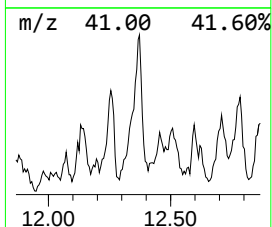
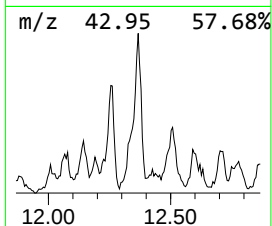
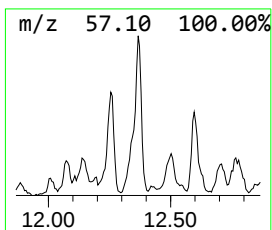
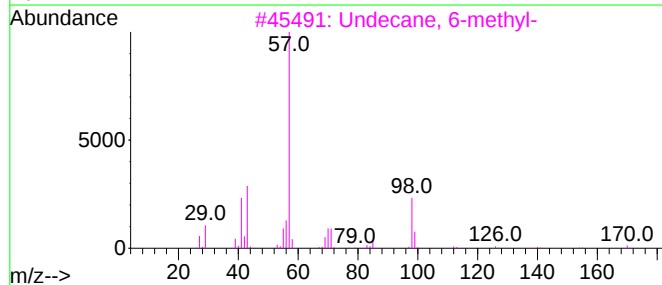
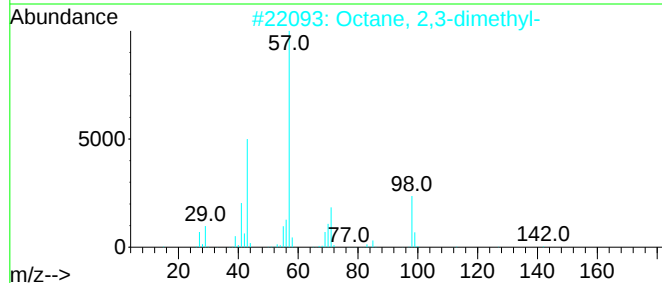
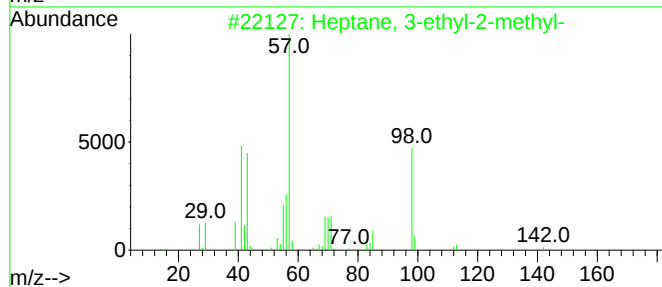
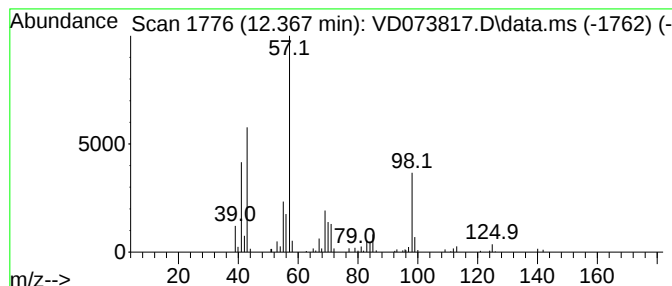
Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D070622S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.367	26.89 ug/l	361121	Chlorobenzene-d5	11.632	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	81
2	Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	72
3	Undecane, 6-methyl-	170	C12H26	017302-33-9	64
4	Heptane, 4-propyl-	142	C10H22	003178-29-8	59
5	Carbonic acid, heptyl vinyl ester	186	C10H18O3	1010383-25-4	50



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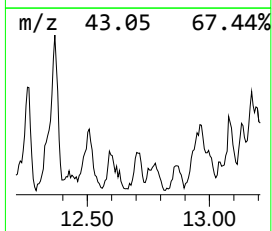
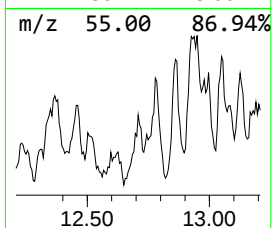
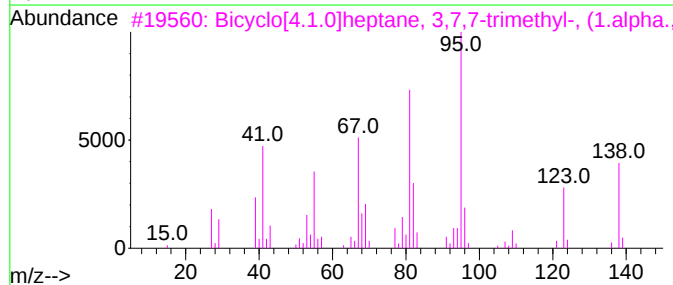
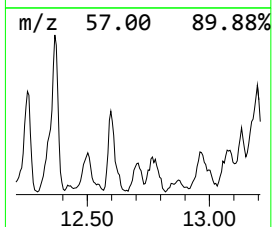
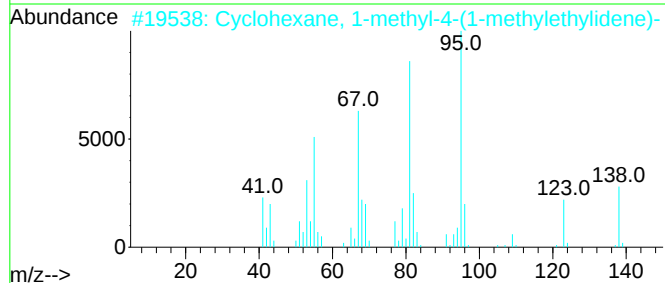
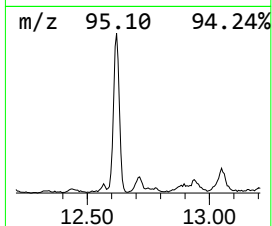
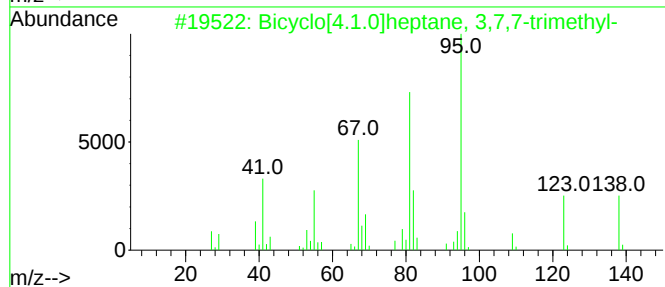
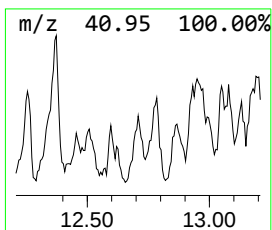
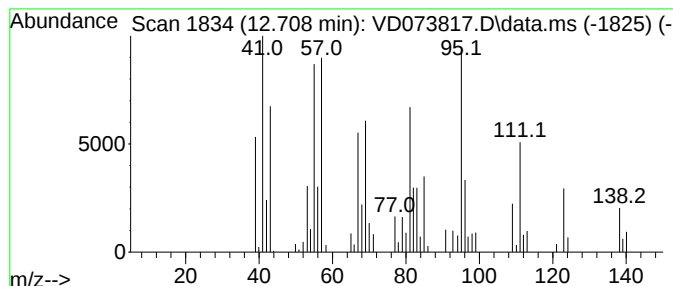
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Quant Title : SW846 8260

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

Peak Number 4 Bicyclo[4.1.0]heptane, 3,7,... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.708	14.50 ug/l	168338	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[4.1.0]heptane, 3,7,7-tri...	138	C10H18	000554-59-6	60
2			Cyclohexane, 1-methyl-4-(1-methy...	138	C10H18	001124-27-2	55
3			Bicyclo[4.1.0]heptane, 3,7,7-tri...	138	C10H18	018968-23-5	55
4			(2E,4E)-3,7-Dimethylocta-2,4-diene	138	C10H18	006874-39-1	46
5			Bicyclo[3.1.0]hexan-2-one, 5-(1-...	138	C9H14O	000513-20-2	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070722\
Data File : VD073817.D
Acq On : 07 Jul 2022 19:28
Operator : VA/SY
Sample : N3593-10
Misc : 4.77G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
20220486A-4A

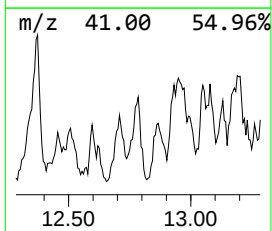
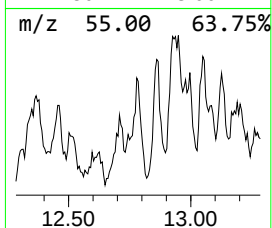
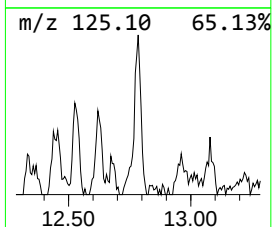
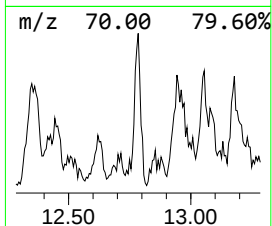
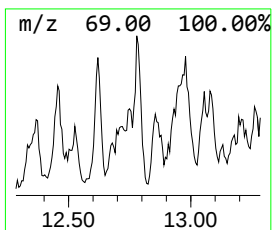
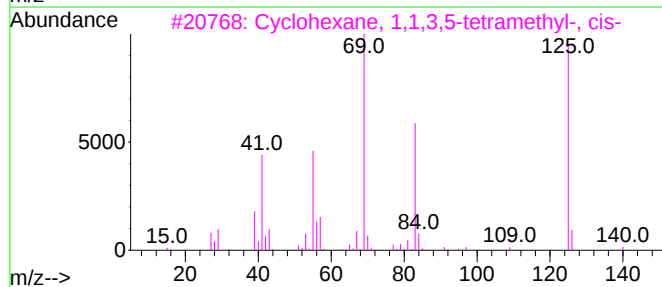
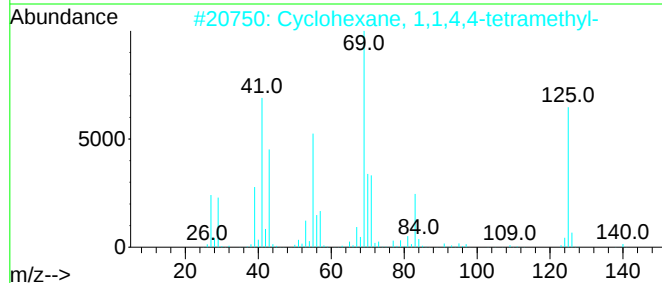
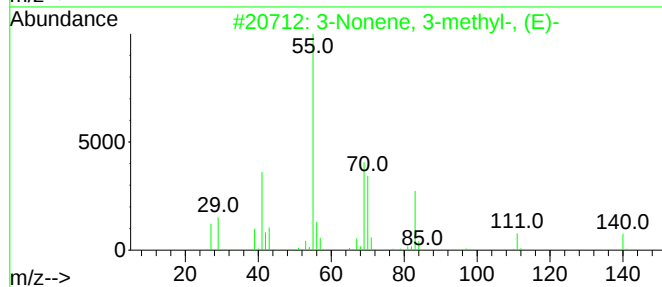
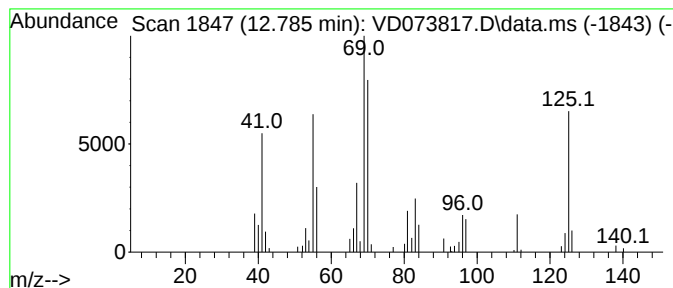
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Quant Title : SW846 8260

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

Peak Number 5 3-Nonene, 3-methyl-, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.785	13.06 ug/l	151627	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Nonene, 3-methyl-, (E)-	140	C10H20	069405-42-1	50
2			Cyclohexane, 1,1,4,4-tetramethyl-	140	C10H20	002223-52-1	49
3			Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-32-9	38
4			2,2-Dimethyl-3-heptene trans	126	C9H18	019550-75-5	35
5			1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070722\
Data File : VD073817.D
Acq On : 07 Jul 2022 19:28
Operator : VA/SY
Sample : N3593-10
Misc : 4.77G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
20220486A-4A

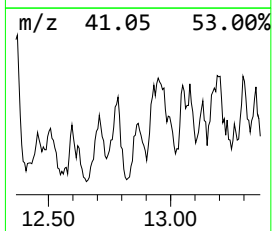
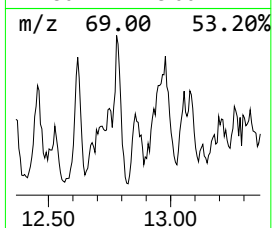
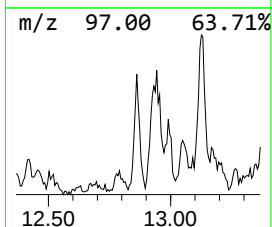
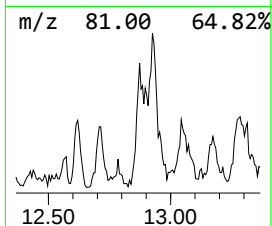
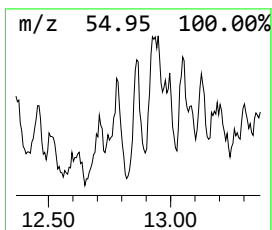
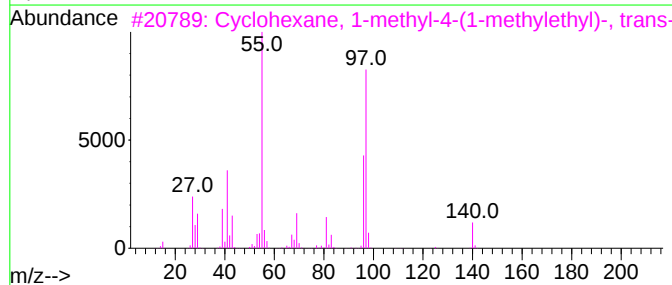
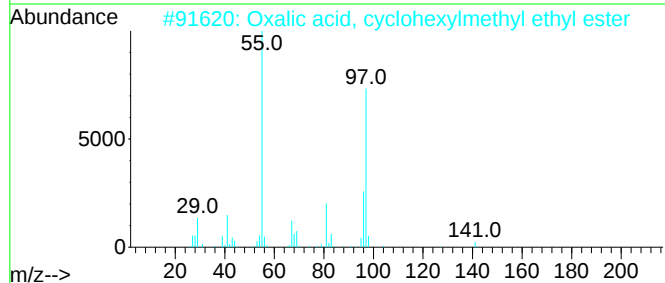
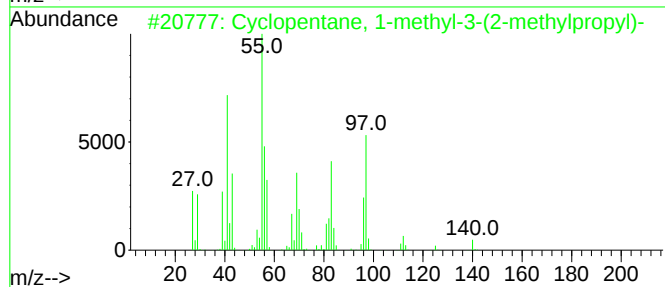
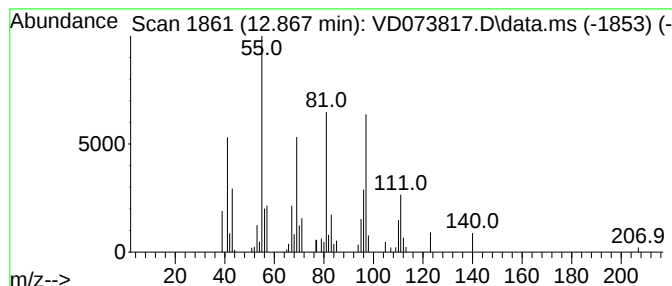
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Quant Title : SW846 8260

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

Peak Number 6 unknown12.867 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.867	13.31 ug/l	154545	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1-methyl-3-(2-meth...	140	C10H20	029053-04-1	49
2			Oxalic acid, cyclohexylmethyl et...	214	C11H18O4	1000309-68-0	47
3			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	46
4			Cyclohexane, 1-methyl-2-propyl-	140	C10H20	004291-79-6	45
5			m-Menthane, (1S,3S)-(+)-	140	C10H20	013837-67-7	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070722\
Data File : VD073817.D
Acq On : 07 Jul 2022 19:28
Operator : VA/SY
Sample : N3593-10
Misc : 4.77G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
20220486A-4A

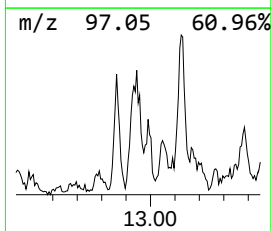
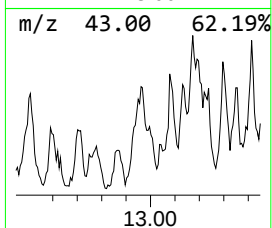
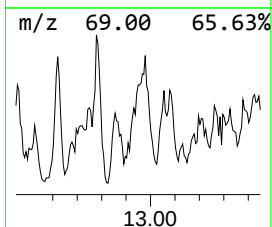
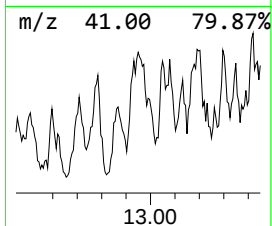
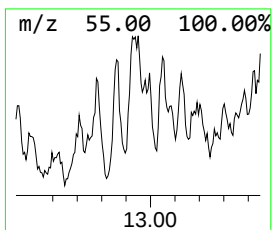
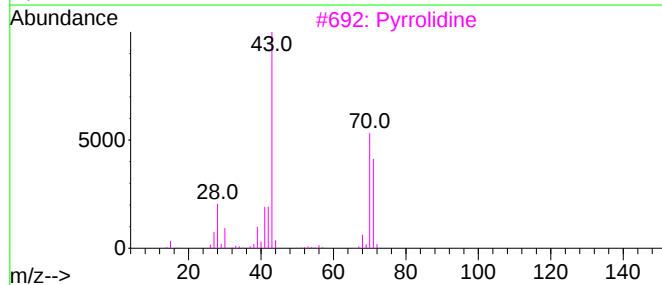
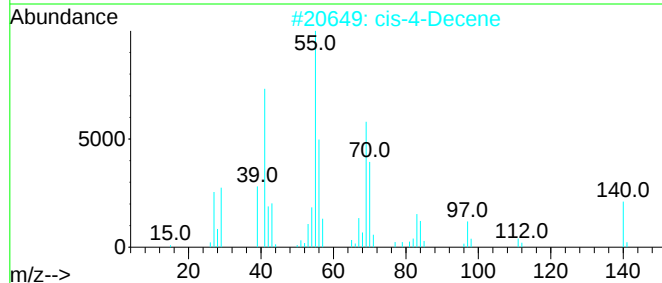
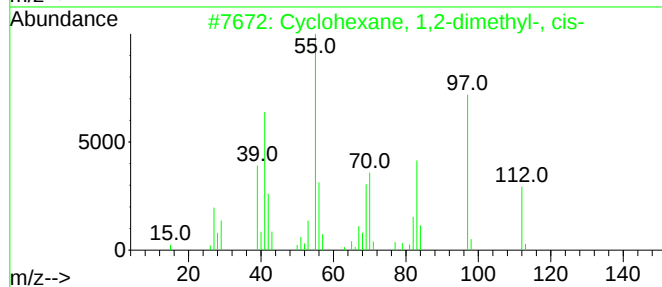
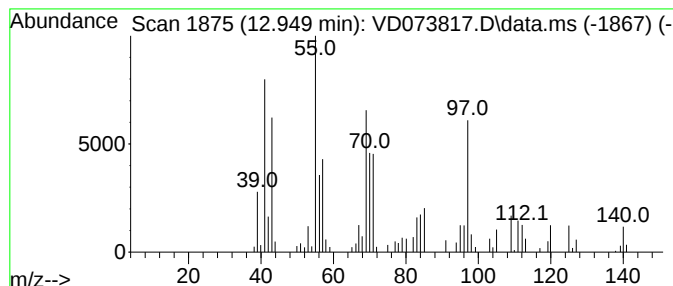
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Quant Title : SW846 8260

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

Peak Number 7 unknown12.949 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.949	26.19 ug/l	304043	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	41
2			cis-4-Decene	140	C10H20	019398-88-0	35
3			Pyrrolidine	71	C4H9N	000123-75-1	30
4			Cyclopropane, 1-ethyl-2-pentyl-	140	C10H20	062238-08-8	30
5			4-Octene, 2,6-dimethyl-, [S-(Z)]-	140	C10H20	062960-77-4	25



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070722\
Data File : VD073817.D
Acq On : 07 Jul 2022 19:28
Operator : VA/SY
Sample : N3593-10
Misc : 4.77G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

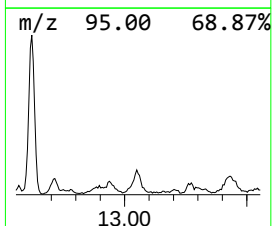
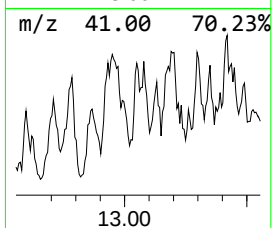
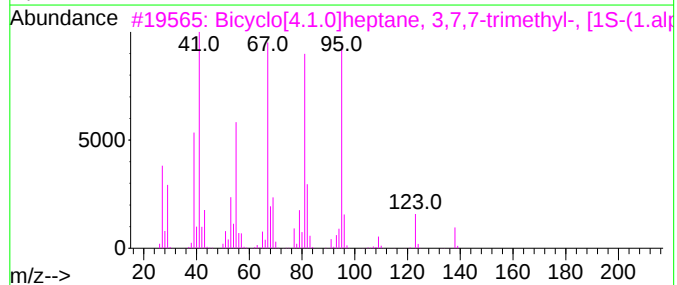
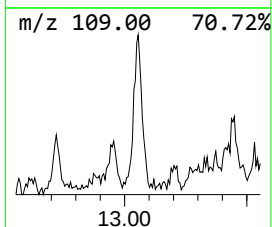
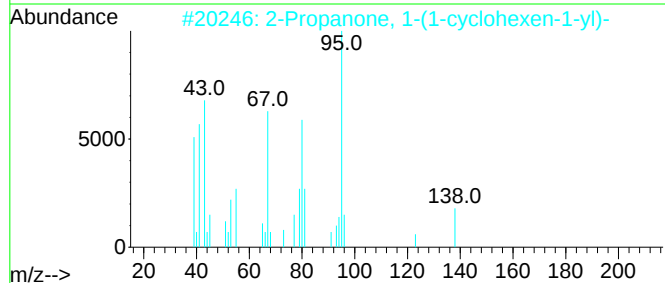
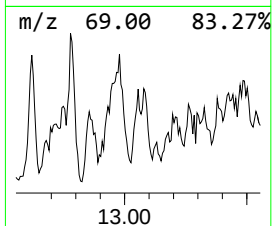
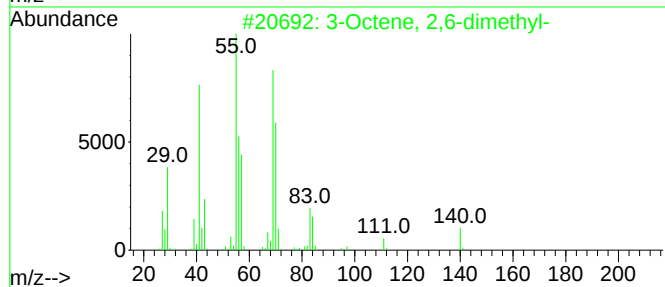
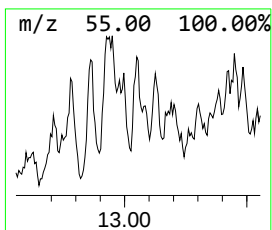
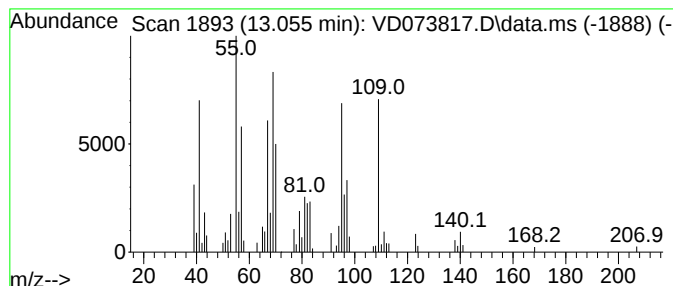
Instrument :
MSVOA_D
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D070622S.M
Quant Title : SW846 8260

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

Peak Number 8 unknown13.055 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD			R.T.	
13.055	10.79 ug/l	125306	1,4-Dichlorobenzene-d4			13.561	
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Octene, 2,6-dimethyl-		140	C10H20		006874-28-8	38
2	2-Propanone, 1-(1-cyclohexen-1-yl)-		138	C9H14O		000768-50-3	35
3	Bicyclo[4.1.0]heptane, 3,7,7-tri...		138	C10H18		006069-97-2	35
4	Cyclopentane, (3-methylbutylidene)-		138	C10H18		053366-51-1	25
5	1-Pentene, 3,3-dimethyl-		98	C7H14		003404-73-7	18



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070722\
Data File : VD073817.D
Acq On : 07 Jul 2022 19:28
Operator : VA/SY
Sample : N3593-10
Misc : 4.77G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

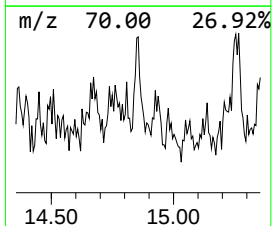
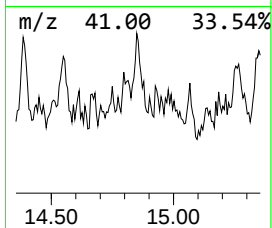
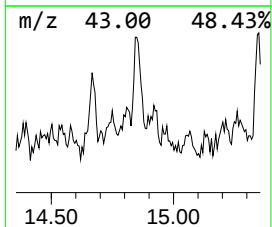
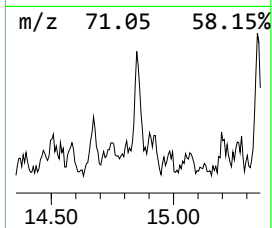
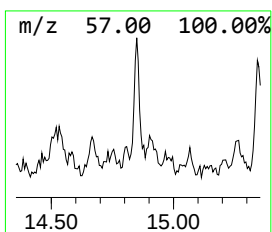
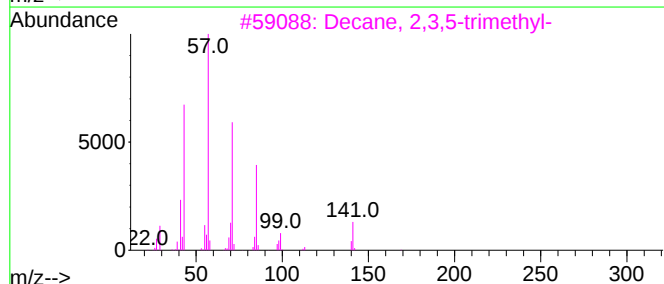
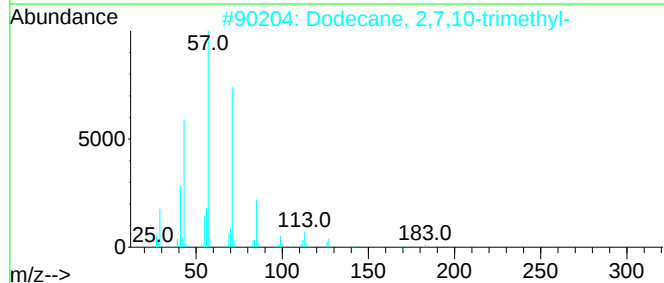
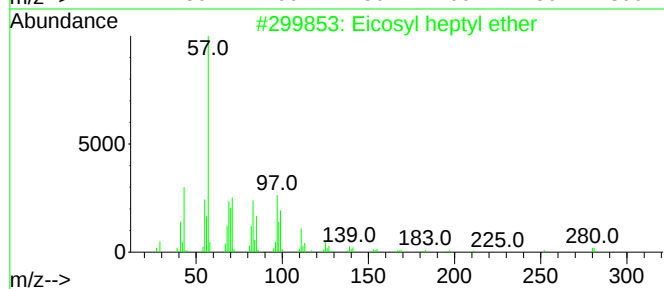
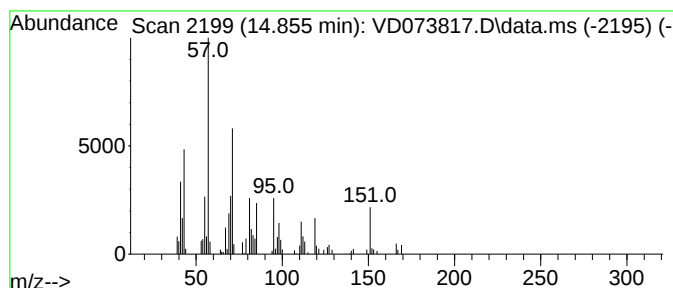
Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D070622S.M
Quant Title : SW846 8260

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

Peak Number 9 unknown14.855 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.855	9.88 ug/l	114756	1,4-Dichlorobenzene-d4	13.561	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosyl heptyl ether	396	C27H56O	1000406-39-6	43
2	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	38
3	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	38
4	Hexadecane	226	C16H34	000544-76-3	38
5	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070722\
Data File : VD073817.D
Acq On : 07 Jul 2022 19:28
Operator : VA/SY
Sample : N3593-10
Misc : 4.77G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
20220486A-4A

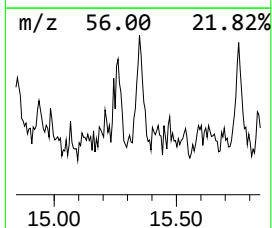
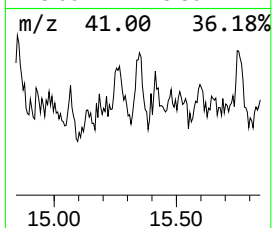
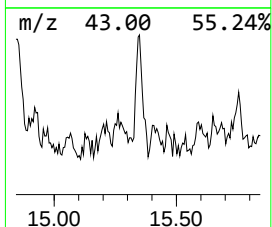
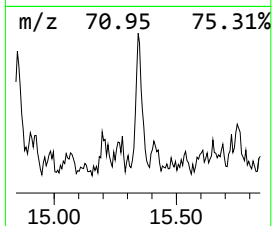
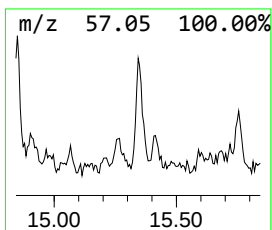
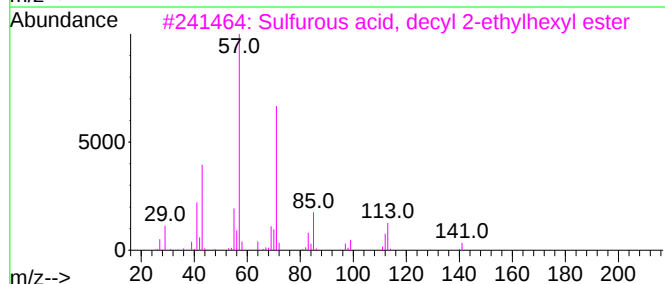
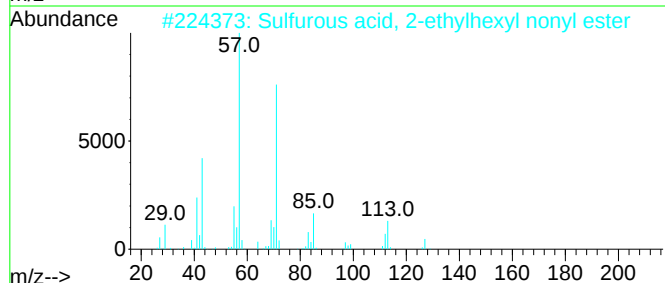
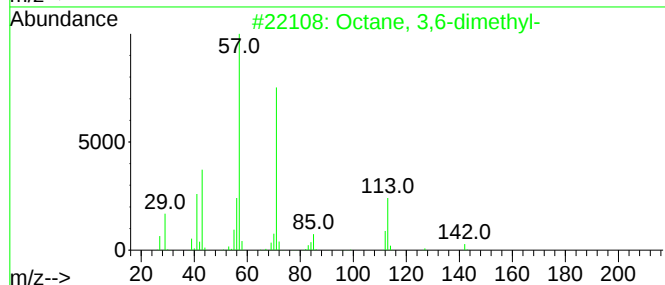
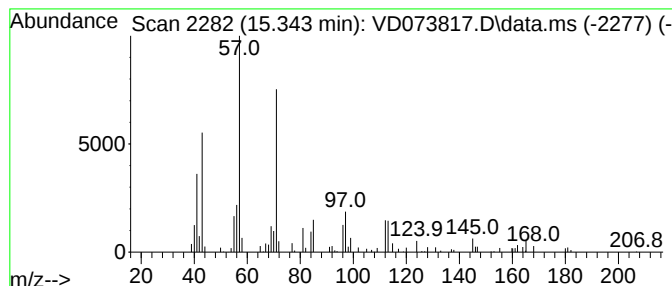
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D070622S.M
Quant Title : SW846 8260

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

Peak Number 10 Octane, 3,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.343	9.14 ug/l	106151	1,4-Dichlorobenzene-d4	13.561

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	59
2			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	59
3			Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	59
4			Oxalic acid, 2-ethylhexyl isobut...	258	C14H26O4	1010309-38-5	53
5			Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070722\
Data File : VD073817.D
Acq On : 07 Jul 2022 19:28
Operator : VA/SY
Sample : N3593-10
Misc : 4.77G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
20220486A-4A

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D070622S.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
Pentane, 2,3,3-...	9.908	9.7	ug/l	103581	2	8.855	532909	50.0
Octane, 2,6-dim...	12.255	11.5	ug/l	154164	3	11.632	671556	50.0
Heptane, 3-ethy...	12.367	26.9	ug/l	361121	3	11.632	671556	50.0
Bicyclo[4.1.0]h...	12.708	14.5	ug/l	168338	4	13.561	580499	50.0
3-Nonene, 3-met...	12.785	13.1	ug/l	151627	4	13.561	580499	50.0
unknown12.867	12.867	13.3	ug/l	154545	4	13.561	580499	50.0
unknown12.949	12.949	26.2	ug/l	304043	4	13.561	580499	50.0
unknown13.055	13.055	10.8	ug/l	125306	4	13.561	580499	50.0
unknown14.855	14.855	9.9	ug/l	114756	4	13.561	580499	50.0
Octane, 3,6-dim...	15.343	9.1	ug/l	106151	4	13.561	580499	50.0