

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD040522\
 Data File : VD072535.D
 Acq On : 05 Apr 2022 15:17
 Operator : VA/SY
 Sample : N2247-09
 Misc : 4.80G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 WC-14A-1A-4-GR-15

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\82D032922S.M
 Title : SW846 8260

Signal : TIC: VD072535.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.038	178	190	209	rBV	3399105	7522002	1.70%	0.132%
2	4.144	363	378	394	rBV	2064073	6915326	1.56%	0.122%
3	4.950	500	515	520	rBV	6439231	22923483	5.17%	0.404%
4	5.014	522	526	560	rVB	7929342	26522303	5.98%	0.467%
5	5.497	589	608	645	rBV	21669910	78585348	17.72%	1.384%
6	6.767	801	824	839	rBV	7163401	24716562	5.57%	0.435%
7	6.932	839	852	861	rVV	11520105	37262895	8.40%	0.656%
8	7.032	861	869	888	rVV	8214798	25100302	5.66%	0.442%
9	7.208	888	899	916	rVB	2571884	8440059	1.90%	0.149%
10	7.738	968	989	1006	rBV	6894865	20730721	4.68%	0.365%
11	7.979	1006	1030	1038	rBV3	33431769	111438987	25.13%	1.962%
12	8.067	1038	1045	1056	rVB	29803812	80074861	18.06%	1.410%
13	8.197	1056	1067	1072	rBV	32316022	84352877	19.02%	1.485%
14	8.255	1072	1077	1086	rVB	23708700	55875057	12.60%	0.984%
15	8.467	1100	1113	1119	rBV2	38841193	104121497	23.48%	1.834%
16	8.538	1119	1125	1130	rVV2	30875753	73708201	16.62%	1.298%
17	8.603	1130	1136	1147	rVB	46899358	115112527	25.96%	2.027%
18	8.861	1170	1180	1186	rBV2	3739558	9360873	2.11%	0.165%
19	9.179	1223	1234	1247	rVB2	9607293	28490724	6.43%	0.502%
20	9.367	1247	1266	1273	rBV4	106485067	443391642	100.00%	7.808%
21	9.538	1289	1295	1300	rVB	12621208	22404956	5.05%	0.395%
22	9.632	1300	1311	1319	rVB2	42932976	118635227	26.76%	2.089%
23	9.773	1319	1335	1345	rBV2	48335760	107719046	24.29%	1.897%
24	9.873	1345	1352	1354	rBV2	6324306	11402131	2.57%	0.201%
25	9.920	1354	1360	1365	rBV	23610932	46903620	10.58%	0.826%
26	9.979	1365	1370	1373	rBV3	16510658	32248654	7.27%	0.568%
27	10.108	1383	1392	1398	rBV4	42856721	136079931	30.69%	2.396%
28	10.232	1406	1413	1417	rBV5	13234424	33437707	7.54%	0.589%
29	10.344	1424	1432	1436	rBV2	86023645	221880052	50.04%	3.907%
30	10.467	1445	1453	1455	rBV2	15042011	26014984	5.87%	0.458%
31	10.514	1455	1461	1471	rVB4	54807197	153507594	34.62%	2.703%
32	10.644	1471	1483	1485	rBV2	27568089	59263571	13.37%	1.044%
33	10.685	1485	1490	1497	rVB2	75297103	165086456	37.23%	2.907%
34	10.779	1497	1506	1511	rBV5	56878779	146370971	33.01%	2.578%
35	10.867	1516	1521	1526	rVB	29986941	54645129	12.32%	0.962%
36	10.961	1526	1537	1543	rBV3	55730145	127907104	28.85%	2.253%

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37	11.061	1543	1554	1562	rBV3	44178506	127151556	28.68%	2.239%
38	11.132	1562	1566	1569	rVV2	31000999	60186116	13.57%	1.060%
39	11.197	1573	1577	1581	rVV2	57308478	120850730	27.26%	2.128%
40	11.249	1581	1586	1592	rVV2	80673529	190344742	42.93%	3.352%
41	11.302	1592	1595	1599	rVV2	27384716	41689460	9.40%	0.734%
42	11.355	1599	1604	1608	rVV3	49091751	96359673	21.73%	1.697%
43	11.449	1608	1620	1628	rVB4	59457771	228954365	51.64%	4.032%
44	11.526	1628	1633	1638	rBV3	42702199	83928092	18.93%	1.478%
45	11.685	1655	1660	1663	rBV3	7163186	11269796	2.54%	0.198%
46	11.738	1664	1669	1672	rBV2	7615717	12783193	2.88%	0.225%
47	11.779	1672	1676	1680	rBV	30333831	49041004	11.06%	0.864%
48	11.826	1680	1684	1690	rVB5	27133498	51534817	11.62%	0.908%
49	11.891	1690	1695	1699	rBV4	50619238	109052405	24.60%	1.920%
50	11.991	1707	1712	1715	rBV2	18414170	33962936	7.66%	0.598%
51	12.079	1719	1727	1732	rVB3	27032729	70014019	15.79%	1.233%
52	12.149	1732	1739	1746	rBV5	74188377	204584126	46.14%	3.603%
53	12.273	1752	1760	1764	rVB3	47173412	104749309	23.62%	1.845%
54	12.355	1764	1774	1776	rBV4	59850348	143300661	32.32%	2.524%
55	12.473	1790	1794	1797	rBV4	12349788	21830413	4.92%	0.384%
56	12.632	1816	1821	1825	rBV3	22630647	40677082	9.17%	0.716%
57	12.714	1829	1835	1839	rBV5	22905784	46483277	10.48%	0.819%
58	12.796	1844	1849	1855	rVB3	71858136	144965444	32.69%	2.553%
59	12.873	1855	1862	1867	rBV6	31208288	76509872	17.26%	1.347%
60	12.955	1868	1876	1881	rBV7	50511728	137677042	31.05%	2.425%
61	13.091	1890	1899	1903	rBV4	31560334	63369292	14.29%	1.116%
62	13.138	1903	1907	1910	rVV3	17877431	32146514	7.25%	0.566%
63	13.179	1910	1914	1922	rVB3	39992491	73681860	16.62%	1.298%
64	13.302	1932	1935	1940	rVB2	11716426	15512268	3.50%	0.273%
65	13.355	1941	1944	1947	rVV	8408381	9992872	2.25%	0.176%
66	13.455	1954	1961	1965	rBV5	33413720	75032917	16.92%	1.321%
67	13.496	1965	1968	1972	rVB3	19481994	26677999	6.02%	0.470%
68	13.596	1982	1985	1989	rVB3	8199170	11887422	2.68%	0.209%
69	13.708	1999	2004	2010	rBV5	18153514	36977178	8.34%	0.651%
70	13.885	2029	2034	2039	rBV2	37531718	66524094	15.00%	1.172%
71	13.926	2039	2041	2046	rVB2	8196030	10934144	2.47%	0.193%
72	13.990	2048	2052	2055	rBV3	7204508	12168634	2.74%	0.214%
73	14.102	2068	2071	2075	rVB	13100118	17939412	4.05%	0.316%
74	14.208	2084	2089	2095	rBV5	12550426	22260374	5.02%	0.392%
75	14.273	2096	2100	2106	rVB6	9412030	16763484	3.78%	0.295%
76	14.332	2107	2110	2112	rBV3	3660632	5180119	1.17%	0.091%

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Integrator: RTE

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Filtering: 5

Sampling : 1

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

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Title : SW846 8260

77	14.396	2116	2121	2126	rBV2	19731041	32984192	7.44%	0.581%
78	14.502	2136	2139	2143	rBV3	3453953	4578834	1.03%	0.081%
79	14.555	2143	2148	2154	rVV6	13823810	23843535	5.38%	0.420%
80	14.808	2186	2191	2195	rVV3	10715425	21041579	4.75%	0.371%
81	14.855	2195	2199	2206	rVB4	11662412	24466253	5.52%	0.431%
82	15.190	2252	2256	2262	rVV4	3562961	6834743	1.54%	0.120%
83	15.261	2265	2268	2278	rVB6	5150461	14054620	3.17%	0.248%
84	15.349	2278	2283	2289	rVB2	6102321	11088957	2.50%	0.195%
85	15.761	2348	2353	2360	rVB6	3235038	6430590	1.45%	0.113%

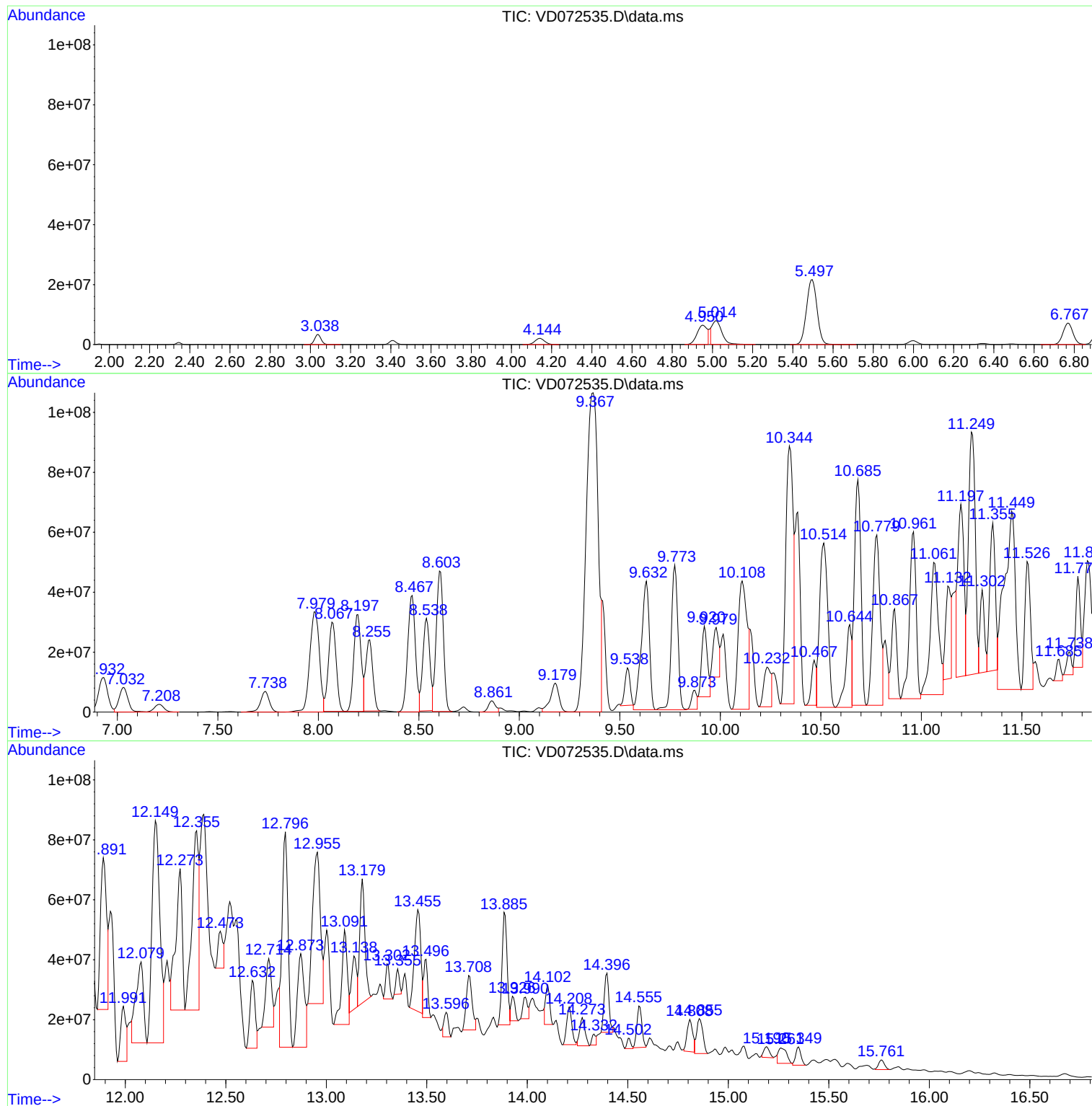
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Instrument :
MSVOA_D
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D032922S.M
Quant Title : SW846 8260

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



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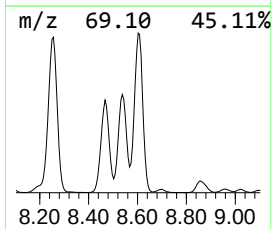
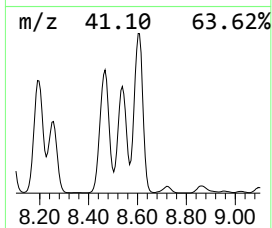
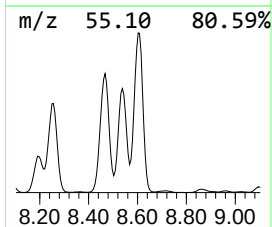
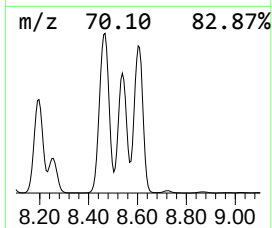
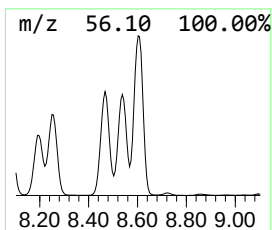
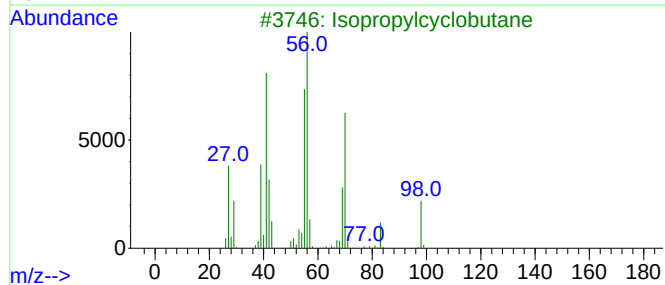
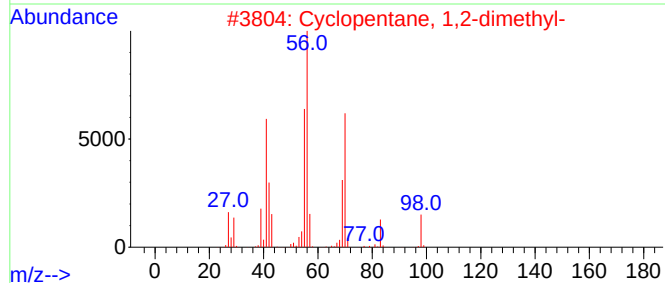
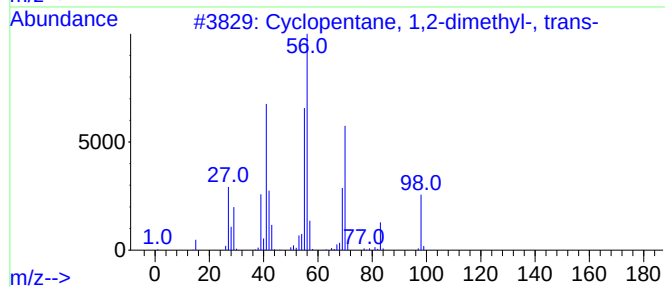
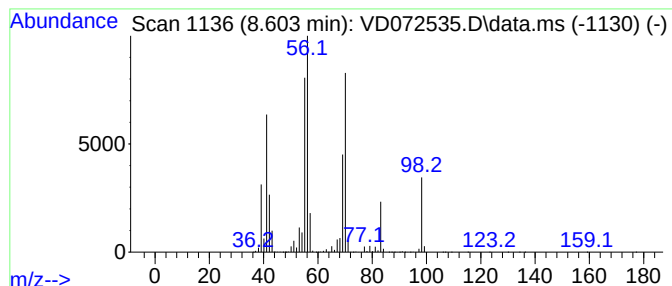
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D032922S.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, 1,2-dimethyl-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.603	614.86 ug/l	115113000	1,4-Difluorobenzene	8.861

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	94
2		Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	94
3		Isopropylcyclobutane	98	C7H14	000872-56-0	94
4		1-Heptene	98	C7H14	000592-76-7	90
5		(Z)-2-Heptene	98	C7H14	006443-92-1	64



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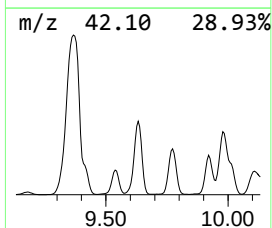
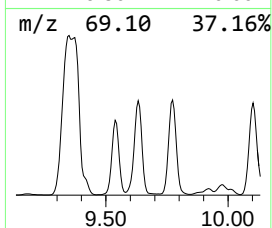
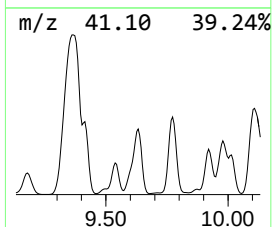
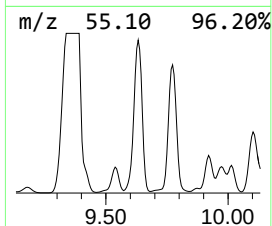
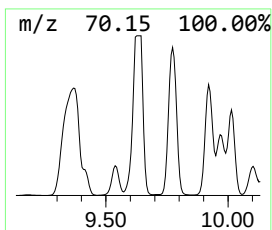
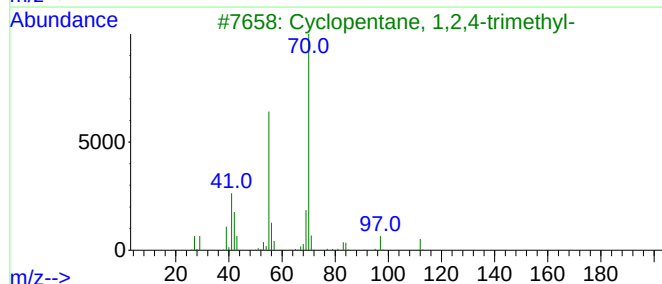
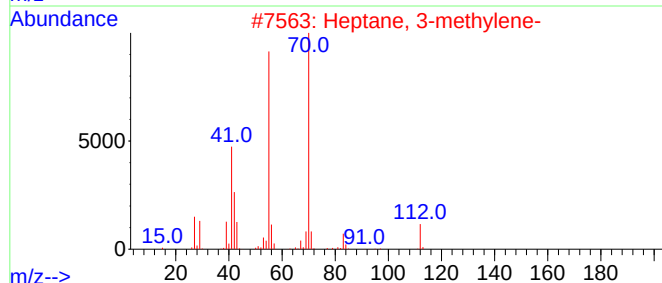
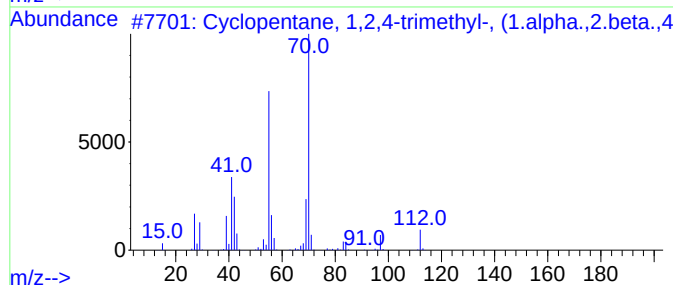
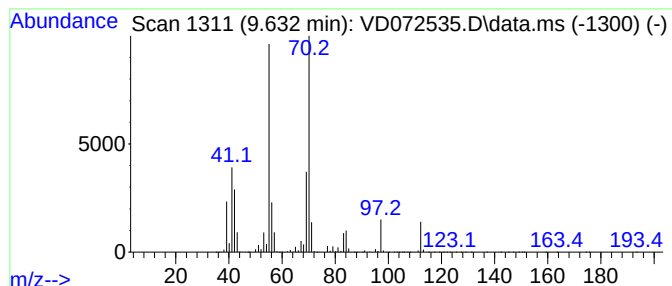
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D032922S.M
Quant Title : SW846 8260

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

Peak Number 2 Cyclopentane, 1,2,4-trimeth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.632	633.68 ug/l	118635000	1,4-Difluorobenzene	8.861

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,2,4-trimethyl-, ...	112	C8H16	016883-48-0	91
2			Heptane, 3-methylene-	112	C8H16	001632-16-2	74
3			Cyclopentane, 1,2,4-trimethyl-	112	C8H16	002815-58-9	53
4			Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	015890-40-1	52
5			Cyclopentane, 1,2,4-trimethyl-, ...	112	C8H16	004850-28-6	30



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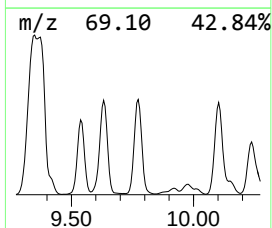
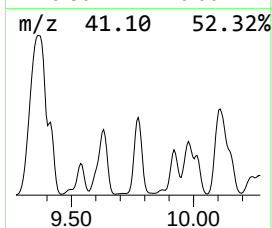
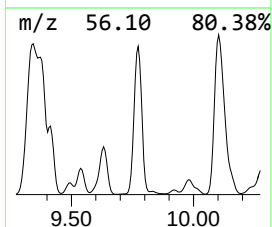
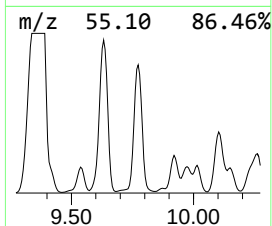
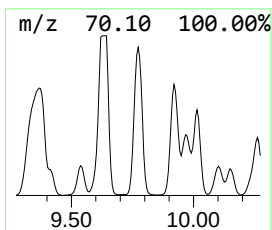
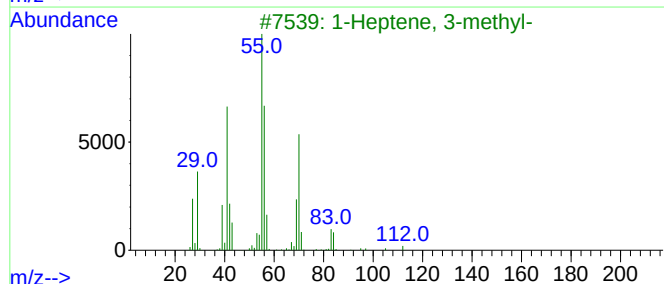
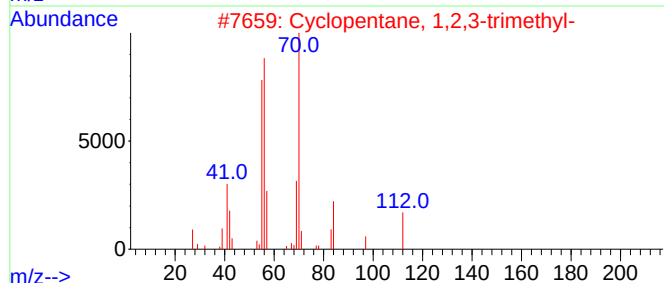
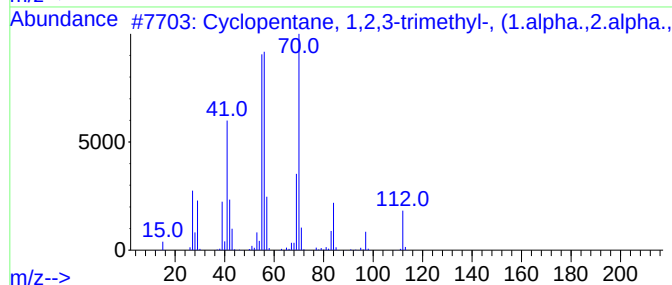
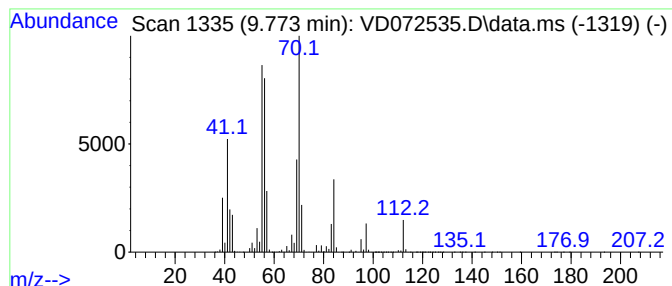
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D032922S.M
Quant Title : SW846 8260

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

Peak Number 3 Cyclopentane, 1,2,3-trimeth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.773	575.37 ug/l	107719000	1,4-Difluorobenzene	8.861

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	015890-40-1	94
2			Cyclopentane, 1,2,3-trimethyl-	112	C8H16	002815-57-8	86
3			1-Heptene, 3-methyl-	112	C8H16	004810-09-7	81
4			Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	59
5			Hexane, 2-methyl-4-methylene-	112	C8H16	003404-80-6	46



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Sample : N2247-09
Misc : 4.80G/5.00ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
WC-14A-1A-4-GR-15

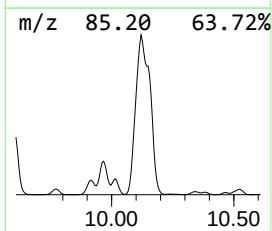
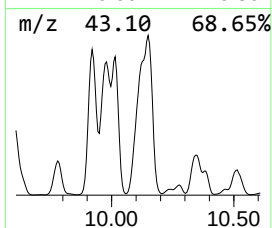
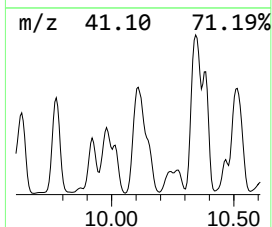
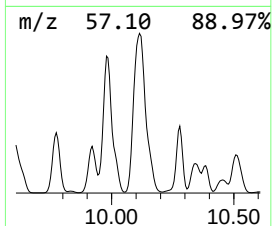
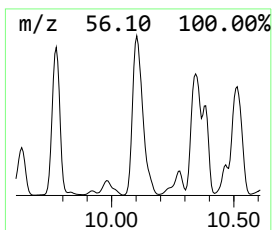
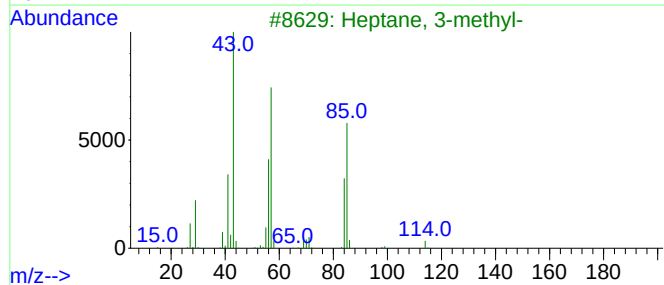
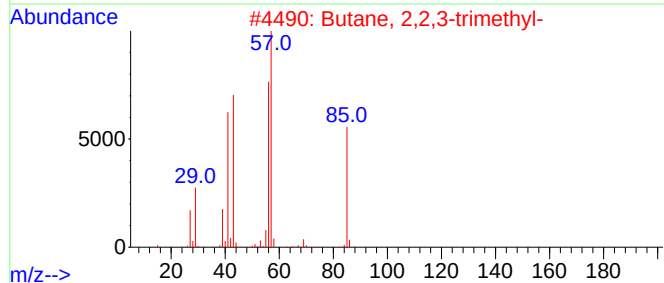
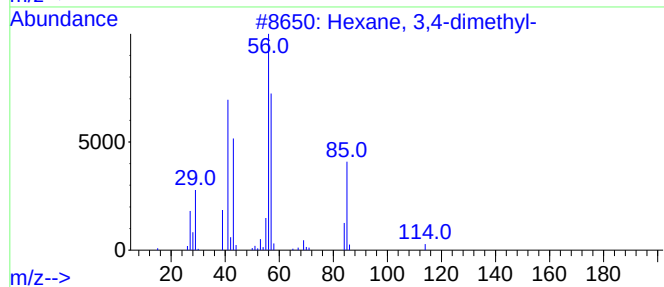
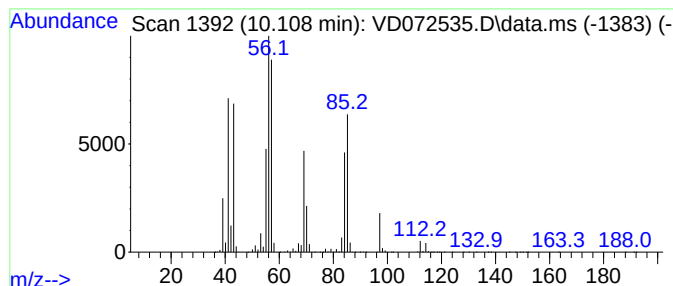
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.108	726.86 ug/l	136080000	1,4-Difluorobenzene	8.861

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3,4-dimethyl-	114	C8H18	000583-48-2	58
2			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	50
3			Heptane, 3-methyl-	114	C8H18	000589-81-1	50
4			Pentanoic acid, butyl ester	158	C9H18O2	000591-68-4	47
5			Butanoic acid, 3-methyl-, butyl ...	158	C9H18O2	000109-19-3	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD040522\
Data File : VD072535.D
Acq On : 05 Apr 2022 15:17
Operator : VA/SY
Sample : N2247-09
Misc : 4.80G/5.00ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
WC-14A-1A-4-GR-15

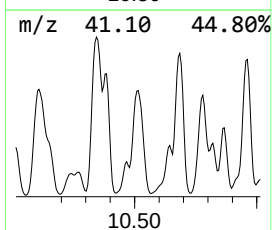
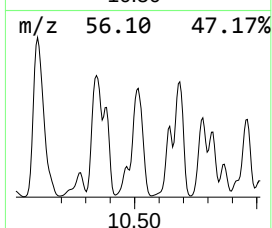
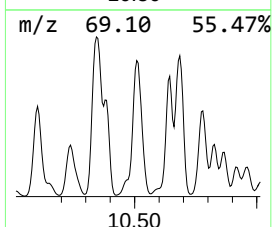
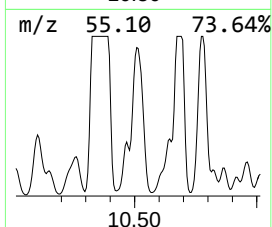
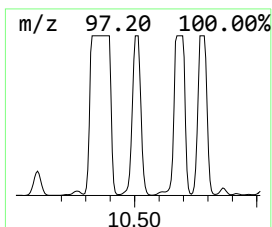
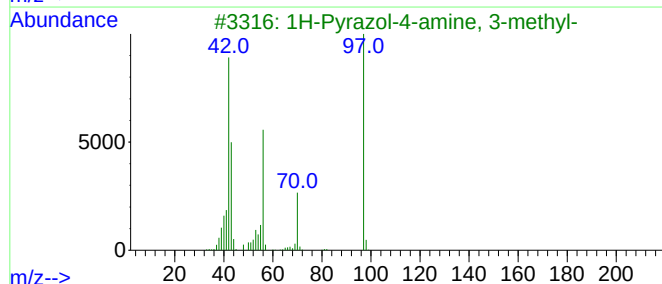
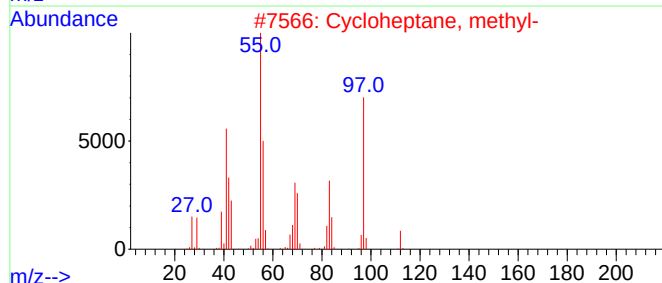
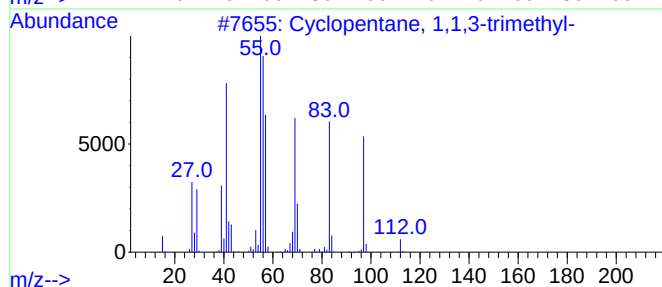
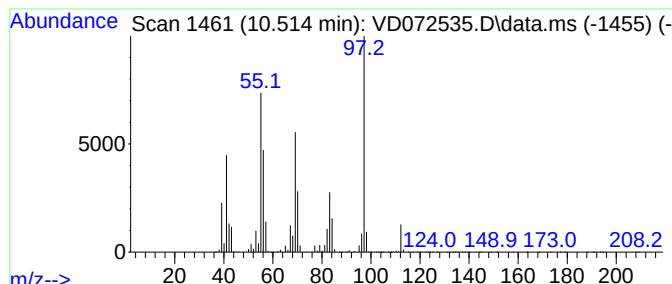
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentane, 1,1,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.514	681.06 ug/l	153508000	Chlorobenzene-d5	11.638

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	72
2			Cycloheptane, methyl-	112	C8H16	004126-78-7	62
3			1H-Pyrazol-4-amine, 3-methyl-	97	C4H7N3	1010338-28-2	52
4			Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	52
5			3-Cyclohexen-1-ol, 3-methyl-	112	C7H12O	053783-91-8	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD040522\
Data File : VD072535.D
Acq On : 05 Apr 2022 15:17
Operator : VA/SY
Sample : N2247-09
Misc : 4.80G/5.00ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
WC-14A-1A-4-GR-15

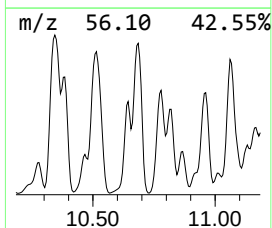
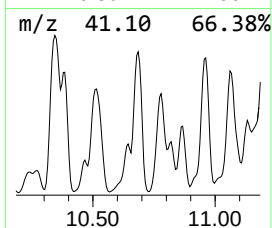
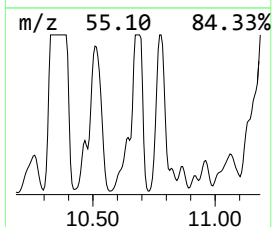
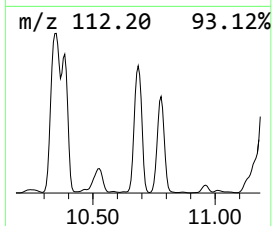
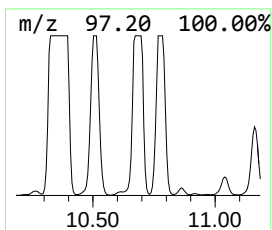
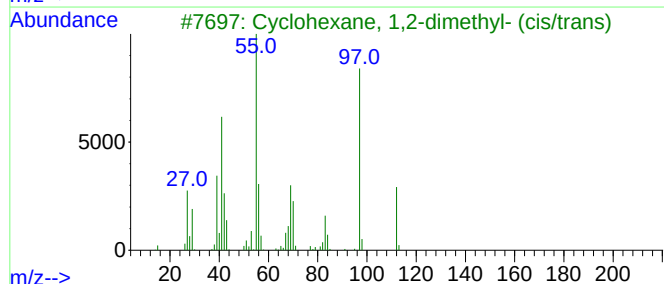
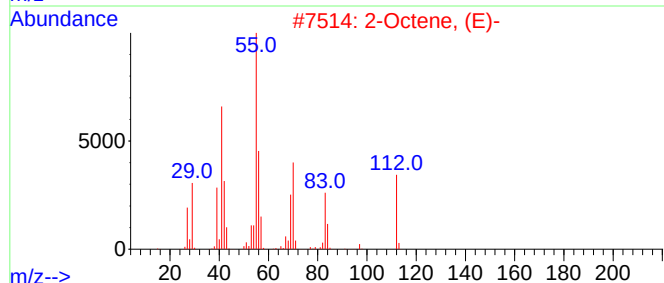
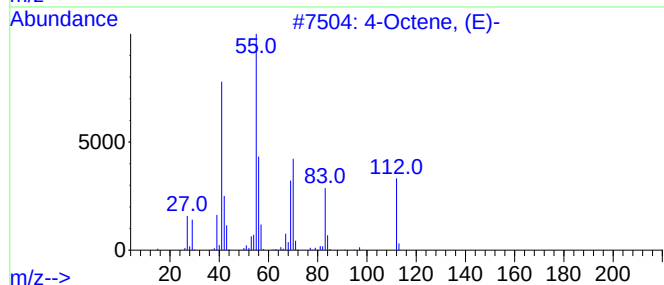
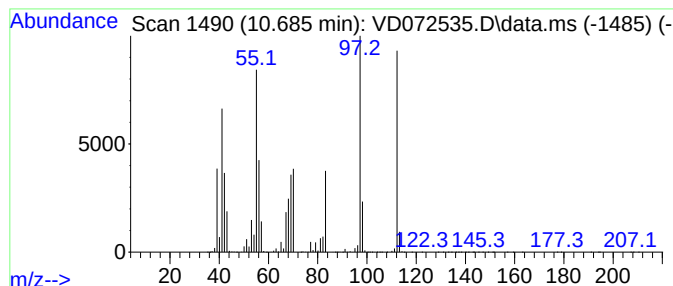
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Quant Title : SW846 8260

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

Peak Number 6 4-Octene, (E)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.685	732.43 ug/l	165086000	Chlorobenzene-d5	11.638

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-Octene, (E)-	112	C8H16	014850-23-8	58
2		2-Octene, (E)-	112	C8H16	013389-42-9	58
3		Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	50
4		3-Octene, (E)-	112	C8H16	014919-01-8	49
5		4-Octene, (Z)-	112	C8H16	007642-15-1	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD040522\
Data File : VD072535.D
Acq On : 05 Apr 2022 15:17
Operator : VA/SY
Sample : N2247-09
Misc : 4.80G/5.00ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
WC-14A-1A-4-GR-15

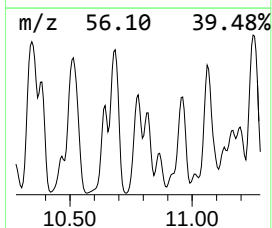
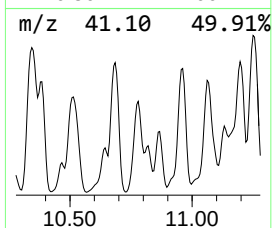
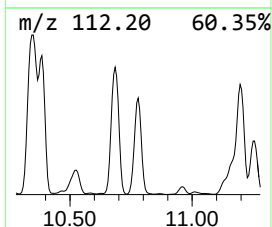
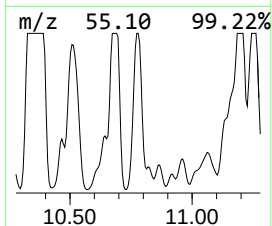
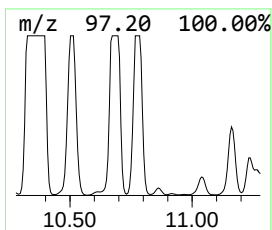
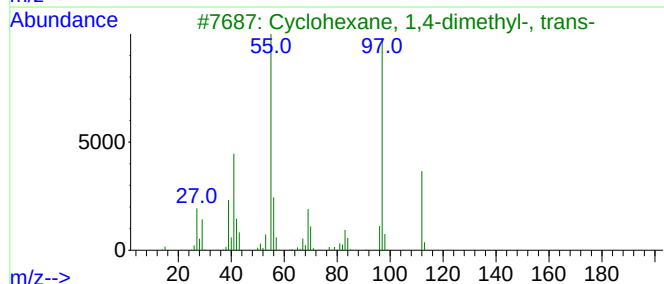
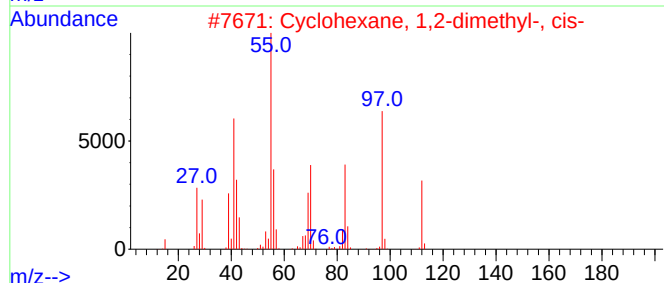
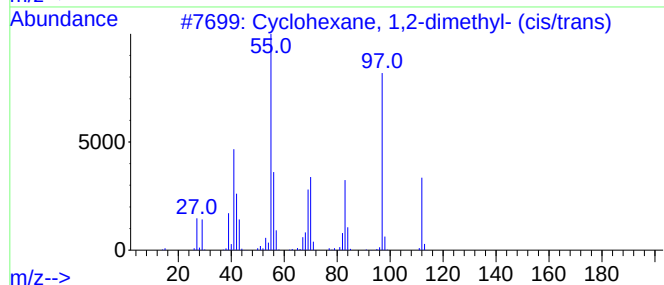
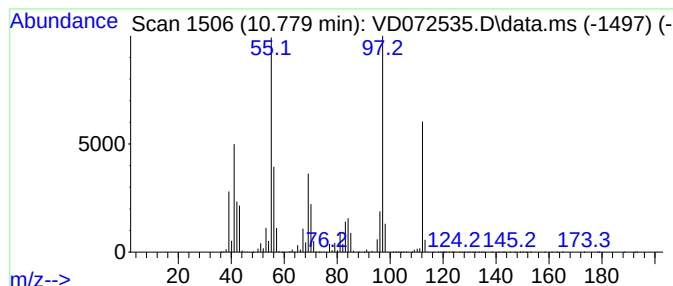
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Quant Title : SW846 8260

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

Peak Number 7 Cyclohexane, 1,2-dimethyl- ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.779	649.39 ug/l	146371000	Chlorobenzene-d5	11.638

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	90
2			Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	87
3			Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	78
4			Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	60
5			Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD040522\
Data File : VD072535.D
Acq On : 05 Apr 2022 15:17
Operator : VA/SY
Sample : N2247-09
Misc : 4.80G/5.00ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

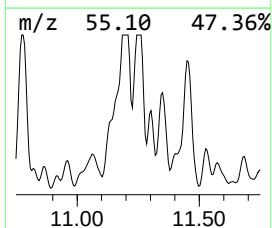
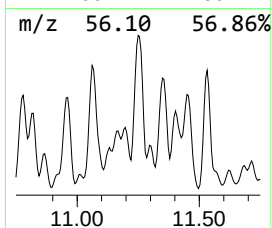
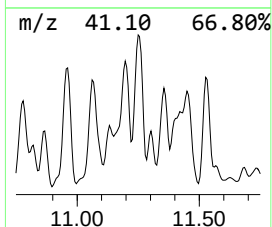
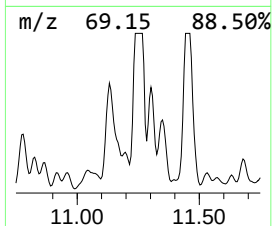
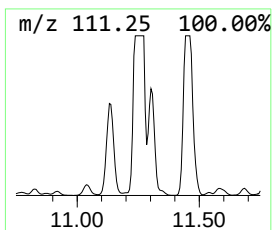
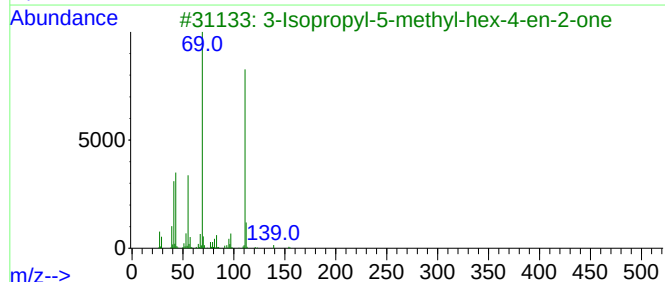
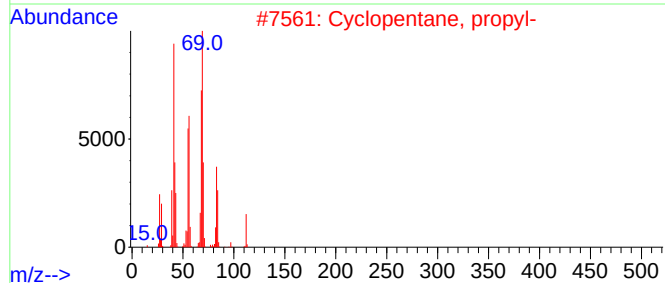
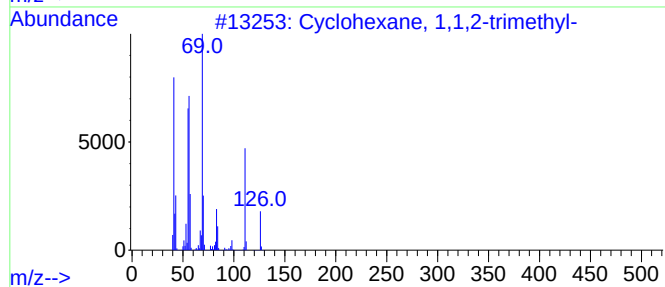
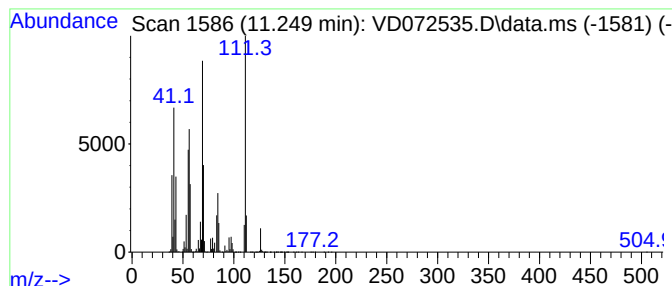
Instrument :
MSVOA_D
ClientSampleId :
WC-14A-1A-4-GR-15

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D032922S.M
Quant Title : SW846 8260

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

Peak Number 8 Cyclohexane, 1,1,2-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.249	844.49 ug/l	190345000	Chlorobenzene-d5	11.638	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	55
2	Cyclopentane, propyl-	112	C8H16	002040-96-2	49
3	3-Isopropyl-5-methyl-hex-4-en-2-one	154	C10H18O	077142-85-9	43
4	4-Methyl-2-heptene	112	C8H16	003404-56-6	41
5	1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD040522\
Data File : VD072535.D
Acq On : 05 Apr 2022 15:17
Operator : VA/SY
Sample : N2247-09
Misc : 4.80G/5.00ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
WC-14A-1A-4-GR-15

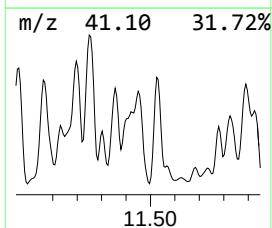
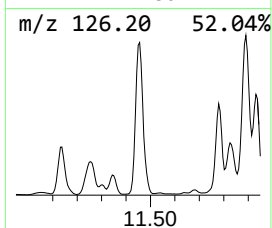
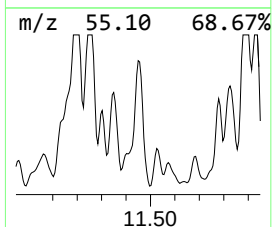
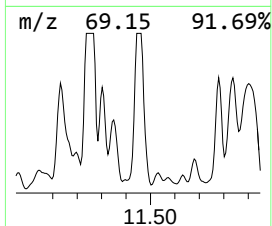
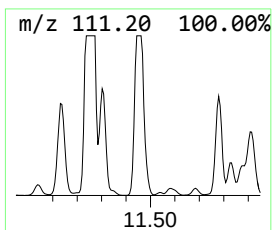
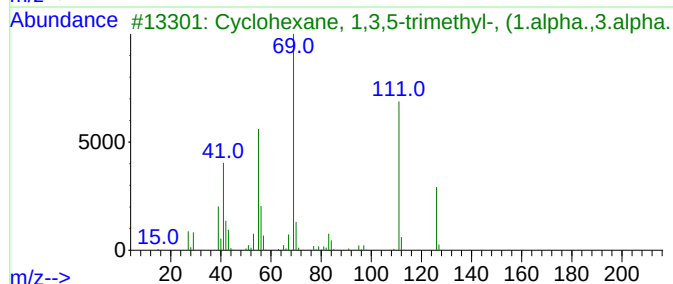
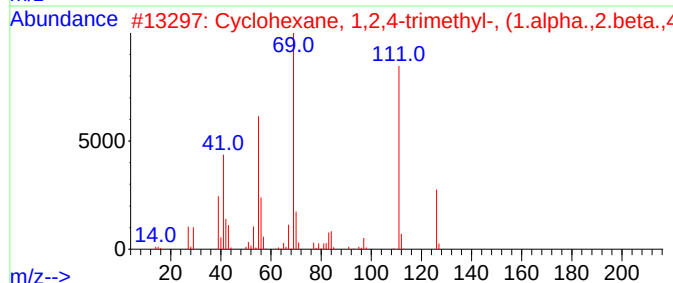
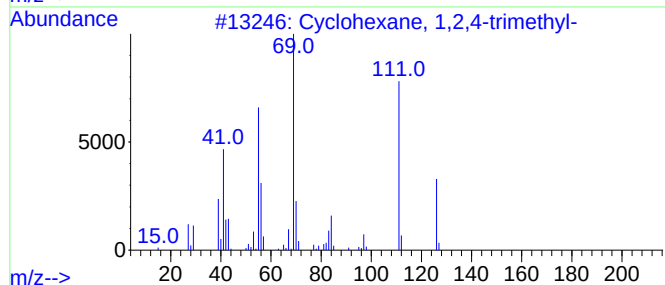
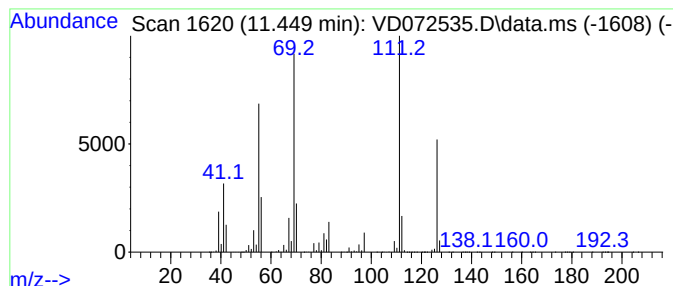
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Quant Title : SW846 8260

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

Peak Number 9 Cyclohexane, 1,2,4-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.449	1015.79 ug/l	228954000	Chlorobenzene-d5	11.638

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	91
2			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	91
3			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	86
4			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	64
5			Cyclooctane, butyl-	168	C12H24	016538-93-5	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD040522\
Data File : VD072535.D
Acq On : 05 Apr 2022 15:17
Operator : VA/SY
Sample : N2247-09
Misc : 4.80G/5.00ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
WC-14A-1A-4-GR-15

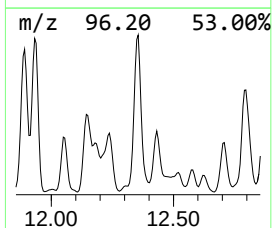
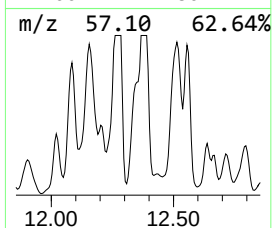
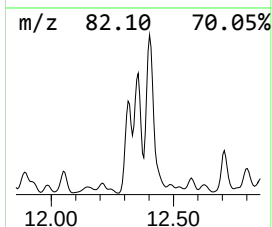
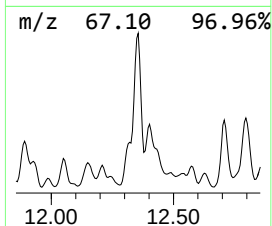
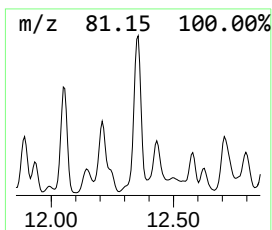
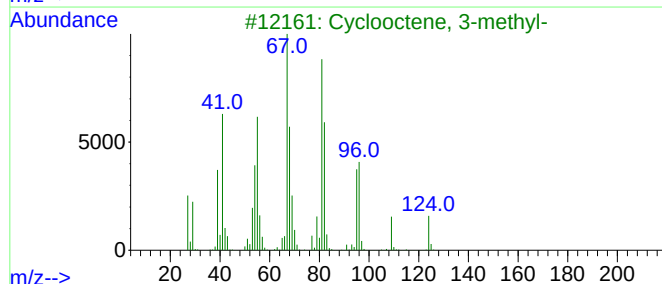
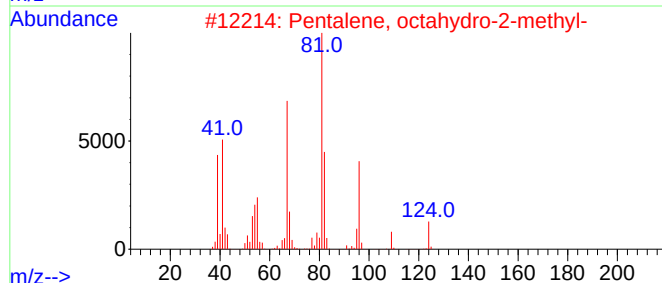
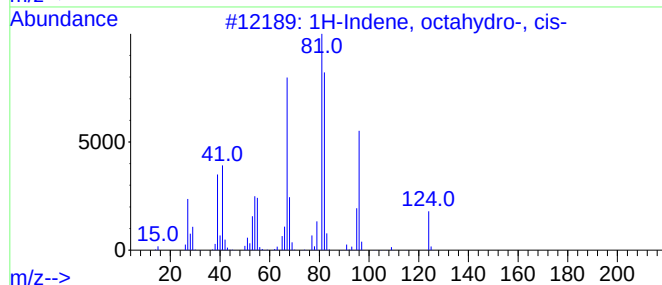
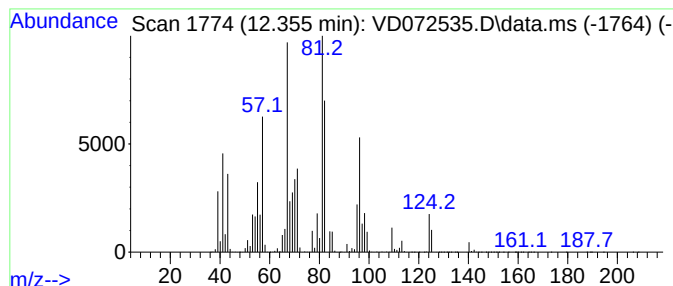
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D032922S.M
Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, octahydro-, cis- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.355	635.77 ug/l	143301000	Chlorobenzene-d5	11.638

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, octahydro-, cis-	124	C9H16	004551-51-3	81
2			Pentalene, octahydro-2-methyl-	124	C9H16	003868-64-2	72
3			Cyclooctene, 3-methyl-	124	C9H16	013152-05-1	70
4			Cyclohexene, 4-propyl-	124	C9H16	013487-64-4	50
5			1H-Indene, octahydro-, trans-	124	C9H16	003296-50-2	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD040522\
Data File : VD072535.D
Acq On : 05 Apr 2022 15:17
Operator : VA/SY
Sample : N2247-09
Misc : 4.80G/5.00ml/MSVOA_D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
WC-14A-1A-4-GR-15

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D032922S.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
Cyclopentane, 1...	8.603	614.9	ug/l	115113000	2	8.861	9360870	50.0
Cyclopentane, 1...	9.632	633.7	ug/l	118635000	2	8.861	9360870	50.0
Cyclopentane, 1...	9.773	575.4	ug/l	107719000	2	8.861	9360870	50.0
Hexane, 3,4-dim...	10.108	726.9	ug/l	136080000	2	8.861	9360870	50.0
Cyclopentane, 1...	10.514	681.1	ug/l	153508000	3	11.638	11269800	50.0
4-Octene, (E)-	10.685	732.4	ug/l	165086000	3	11.638	11269800	50.0
Cyclohexane, 1,...	10.779	649.4	ug/l	146371000	3	11.638	11269800	50.0
Cyclohexane, 1,...	11.249	844.5	ug/l	190345000	3	11.638	11269800	50.0
Cyclohexane, 1,...	11.449	1015.8	ug/l	228954000	3	11.638	11269800	50.0
1H-Indene, octa...	12.355	635.8	ug/l	143301000	3	11.638	11269800	50.0