

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD122121\
 Data File : VD071417.D
 Acq On : 21 Dec 2021 18:16
 Operator : VA/SY
 Sample : M5182-02
 Misc : 7.69G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 BLIND-DUPLICATE

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\82D112921S.M
 Title : SW846 8260

Signal : TIC: VD071417.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.979	1024	1030	1038	rVB2	554045	1227117	1.40%	0.126%
2	8.861	1172	1180	1192	rBV	878716	1796111	2.05%	0.185%
3	9.626	1297	1310	1319	rVB2	563127	1408165	1.60%	0.145%
4	9.767	1325	1334	1347	rVB2	601639	1232588	1.40%	0.127%
5	10.091	1380	1389	1395	rBV	645134	1322861	1.51%	0.136%
6	10.332	1423	1430	1444	rBV2	1545179	3464366	3.95%	0.356%
7	10.508	1454	1460	1470	rVB3	567194	1520236	1.73%	0.156%
8	10.632	1470	1481	1483	rBV2	504681	1056122	1.20%	0.109%
9	10.673	1483	1488	1498	rVB	2639350	4950720	5.64%	0.509%
10	10.785	1498	1507	1510	rBV3	418104	937412	1.07%	0.096%
11	10.949	1525	1535	1542	rVB	1788833	3433326	3.91%	0.353%
12	11.049	1542	1552	1560	rBV	2146856	5022607	5.72%	0.517%
13	11.120	1560	1564	1567	rBV	604437	906880	1.03%	0.093%
14	11.238	1577	1584	1590	rBV2	2440243	4998875	5.69%	0.514%
15	11.297	1590	1594	1597	rVV	1230282	1951946	2.22%	0.201%
16	11.344	1597	1602	1606	rVB	1042363	1770316	2.02%	0.182%
17	11.391	1606	1610	1614	rVB	665409	969722	1.10%	0.100%
18	11.444	1614	1619	1629	rVB	3597061	6558384	7.47%	0.675%
19	11.567	1629	1640	1643	rBV4	591279	1506542	1.72%	0.155%
20	11.638	1646	1652	1656	rBV2	1031500	1747601	1.99%	0.180%
21	11.732	1662	1668	1671	rBV2	605387	980397	1.12%	0.101%
22	11.773	1671	1675	1679	rVB	1892647	2669380	3.04%	0.275%
23	11.826	1679	1684	1689	rBV2	1675268	2991416	3.41%	0.308%
24	11.920	1689	1700	1706	rVB7	3152589	11566698	13.17%	1.190%
25	11.985	1706	1711	1715	rBV2	1314407	2698170	3.07%	0.278%
26	12.079	1724	1727	1731	rVB	982206	1143999	1.30%	0.118%
27	12.149	1731	1739	1744	rBV3	7480617	17362151	19.77%	1.787%
28	12.196	1744	1747	1750	rBV2	1193126	1377398	1.57%	0.142%
29	12.232	1750	1753	1757	rBV3	998052	1455867	1.66%	0.150%
30	12.314	1762	1767	1768	rBV2	3568346	5433430	6.19%	0.559%
31	12.349	1768	1773	1782	rVB4	10003722	20972333	23.89%	2.158%
32	12.432	1783	1787	1789	rBV	1856268	3127338	3.56%	0.322%
33	12.467	1789	1793	1798	rVB3	4445210	7721921	8.79%	0.795%
34	12.514	1798	1801	1808	rVB5	2110715	4187290	4.77%	0.431%
35	12.573	1808	1811	1815	rVB	1599010	2101658	2.39%	0.216%
36	12.626	1815	1820	1825	rVB2	3672602	6468509	7.37%	0.666%

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Integrator: RTE

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
 Title : SW846 8260

37	12.714	1825	1835	1839	rBV5	5719880	15643670	17.82%	1.610%
38	12.791	1843	1848	1855	rVB3	15419093	30043718	34.22%	3.092%
39	12.873	1855	1862	1866	rBV4	7257883	17317042	19.72%	1.782%
40	12.943	1868	1874	1881	rBV6	8369162	25177083	28.67%	2.591%
41	13.055	1890	1893	1894	rBV2	1193543	1402182	1.60%	0.144%
42	13.091	1895	1899	1903	rVV2	7564490	11305668	12.88%	1.163%
43	13.132	1903	1906	1911	rVV3	4643288	8581902	9.77%	0.883%
44	13.196	1911	1917	1927	rVB8	6856150	21561302	24.56%	2.219%
45	13.308	1927	1936	1940	rBV3	11525835	26464415	30.14%	2.723%
46	13.361	1941	1945	1946	rVV3	8686015	12925642	14.72%	1.330%
47	13.391	1947	1950	1954	rVV	14324775	22111853	25.18%	2.275%
48	13.455	1954	1961	1965	rVV5	11201258	30946206	35.24%	3.184%
49	13.496	1965	1968	1973	rVB4	7924547	11755175	13.39%	1.210%
50	13.714	1999	2005	2016	rVB7	13782777	39383783	44.85%	4.053%
51	13.890	2029	2035	2040	rBV2	42173643	73749823	83.99%	7.589%
52	13.943	2040	2044	2048	rVV5	6011092	9974793	11.36%	1.026%
53	13.990	2049	2052	2056	rVV4	9161408	13781128	15.70%	1.418%
54	14.102	2069	2071	2084	rVB3	22872526	47622467	54.24%	4.901%
55	14.208	2084	2089	2095	rBV5	12697586	24884208	28.34%	2.561%
56	14.279	2096	2101	2107	rVB4	17264032	29588855	33.70%	3.045%
57	14.332	2107	2110	2115	rBV3	7932430	12776158	14.55%	1.315%
58	14.396	2115	2121	2125	rBV2	49643712	87805153	100.00%	9.035%
59	14.508	2136	2140	2143	rBV2	8814768	13028042	14.84%	1.341%
60	14.561	2144	2149	2155	rVV5	32261251	58820984	66.99%	6.053%
61	14.749	2178	2181	2186	rBV3	8612363	13146050	14.97%	1.353%
62	14.814	2187	2192	2197	rVV5	16839946	36131365	41.15%	3.718%
63	14.867	2197	2201	2208	rVB5	25362078	46640657	53.12%	4.800%
64	14.990	2219	2222	2225	rBV	4931428	6017001	6.85%	0.619%
65	15.190	2252	2256	2262	rVB3	9636043	18633008	21.22%	1.917%
66	15.349	2279	2283	2290	rVB2	11530266	20440861	23.28%	2.103%
67	15.437	2290	2298	2305	rBV7	9740050	29563094	33.67%	3.042%
68	15.761	2348	2353	2361	rVB6	4840956	10913053	12.43%	1.123%
69	16.408	2459	2463	2469	rVB2	979800	1953797	2.23%	0.201%
70	16.679	2502	2509	2518	rVB7	584868	1650833	1.88%	0.170%

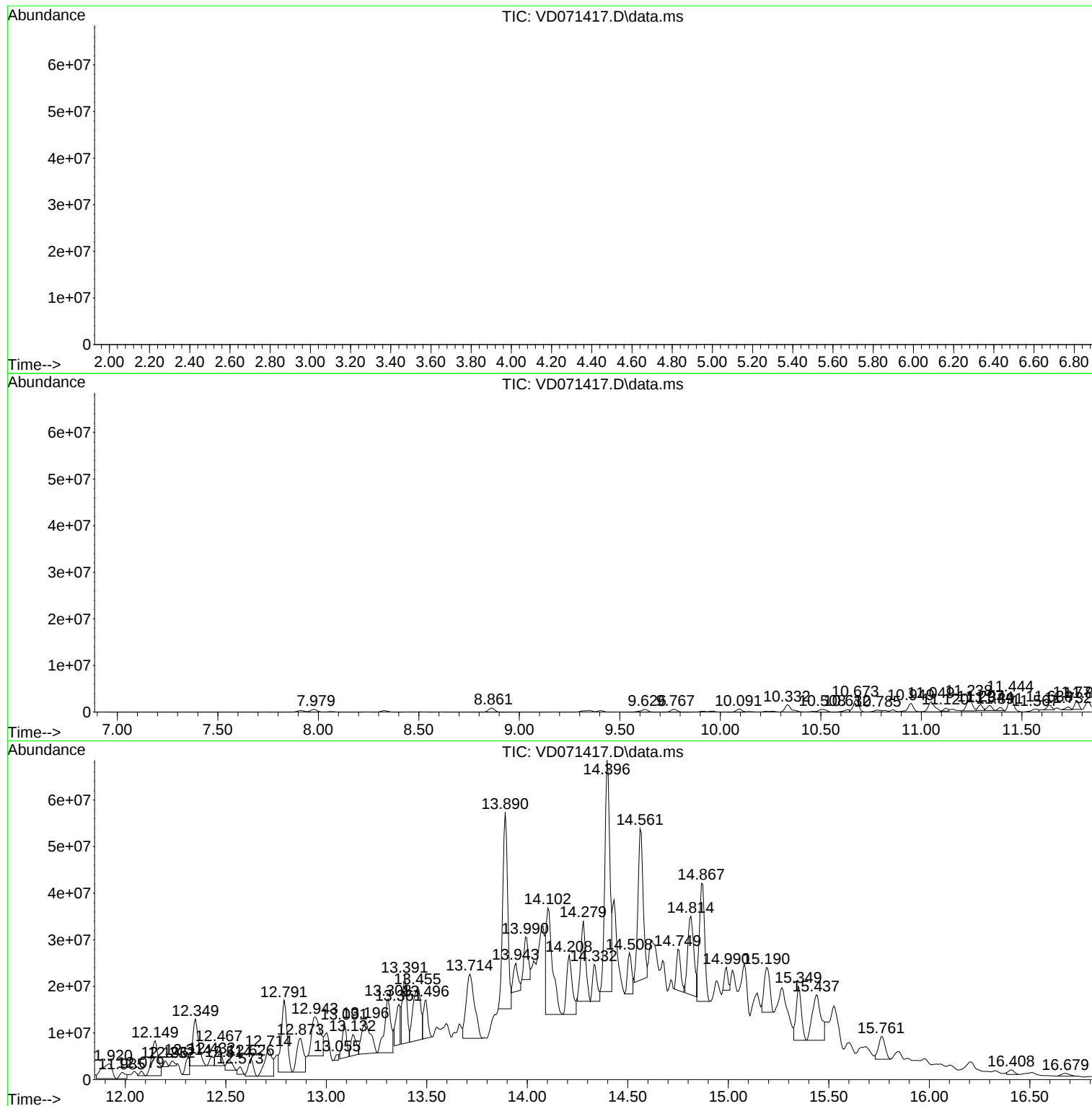
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Instrument :
MSVOA_D
ClientSampleId :
BLIND-DUPLICATE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

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



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Operator : VA/SY
Sample : M5182-02
Misc : 7.69G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

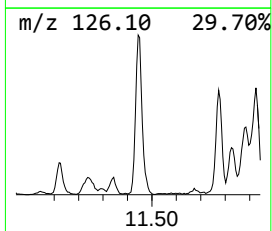
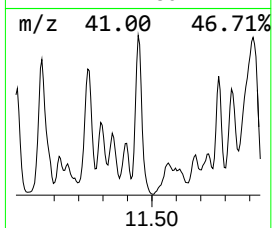
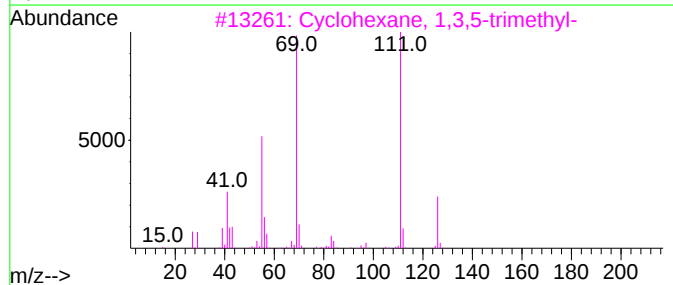
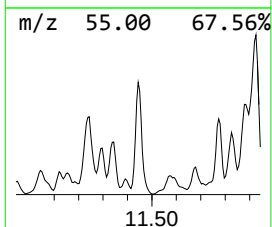
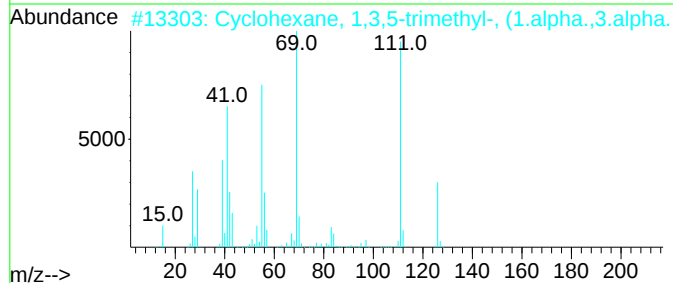
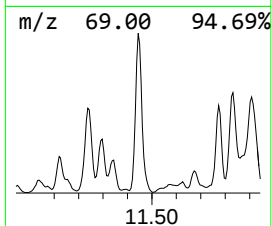
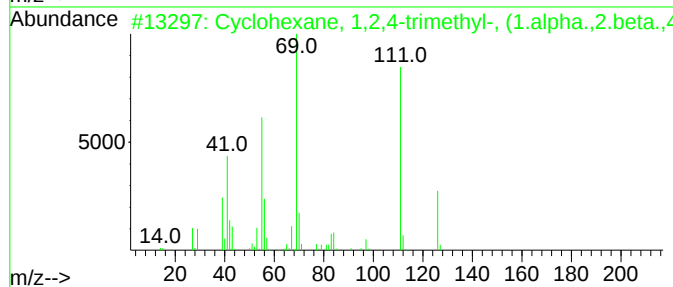
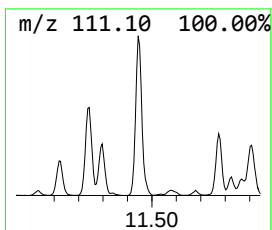
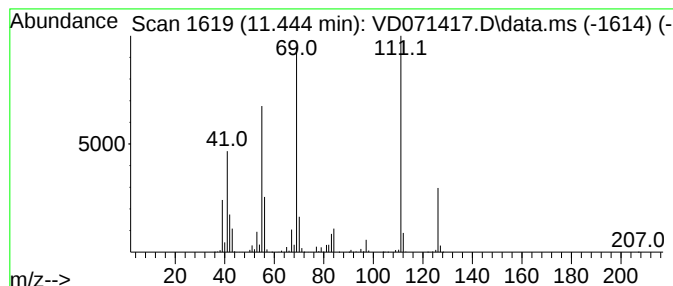
Instrument :
MSVOA_D
ClientSampleId :
BLIND-DUPLICATE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD			R.T.
11.444	187.64 ug/l	6558380	Chlorobenzene-d5			11.638
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2,4-trimethyl-, (...		126	C9H18	007667-60-9	96
2	Cyclohexane, 1,3,5-trimethyl-, (...		126	C9H18	001795-26-2	93
3	Cyclohexane, 1,3,5-trimethyl-		126	C9H18	001839-63-0	87
4	Cyclohexane, 1,3,5-trimethyl-, (...		126	C9H18	001795-27-3	87
5	Cyclohexane, 1,2,4-trimethyl-		126	C9H18	002234-75-5	74



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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BLIND-DUPLICATE

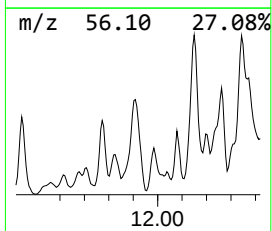
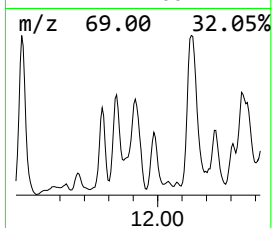
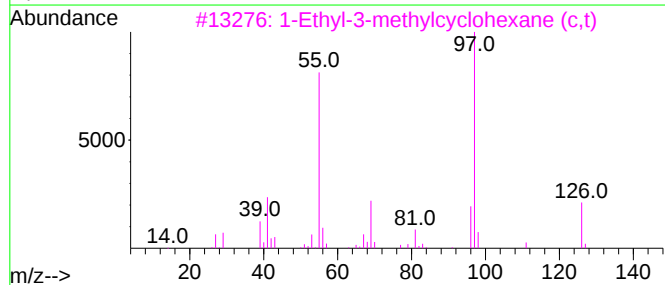
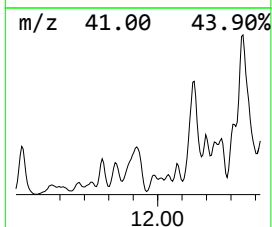
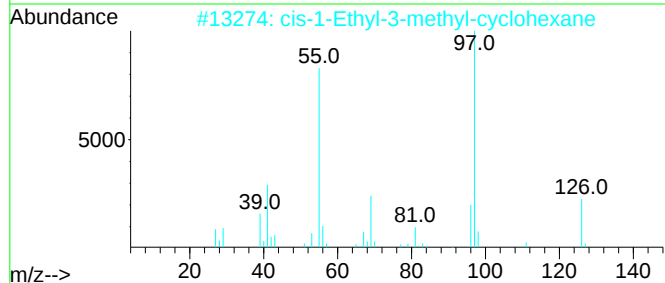
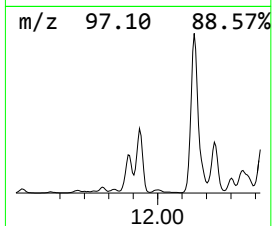
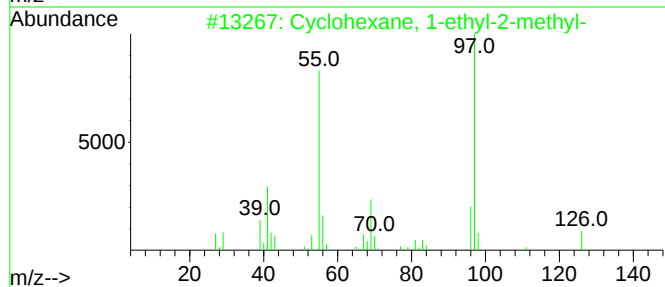
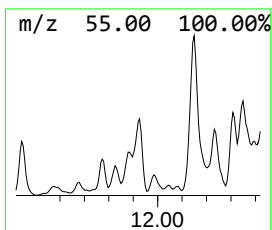
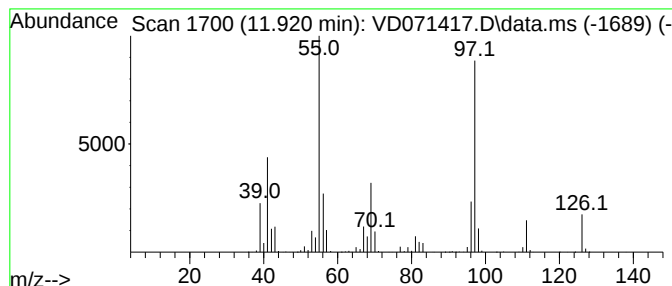
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

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

Peak Number 2 Cyclohexane, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.920	330.93 ug/l	11566700	Chlorobenzene-d5	11.638

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	90
2			cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	81
3			1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	81
4			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	74
5			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	74



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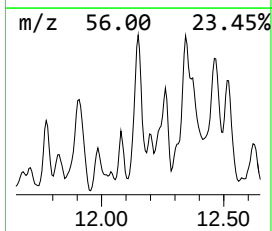
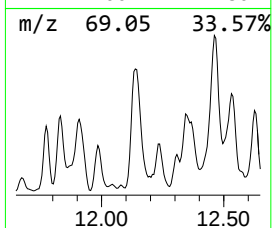
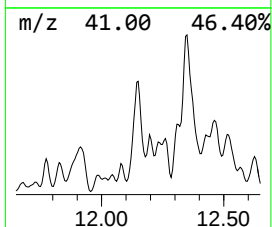
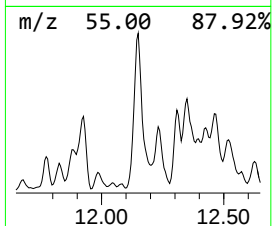
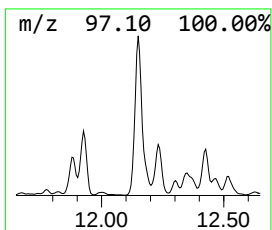
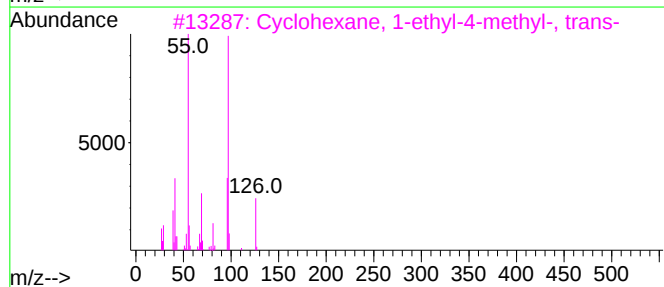
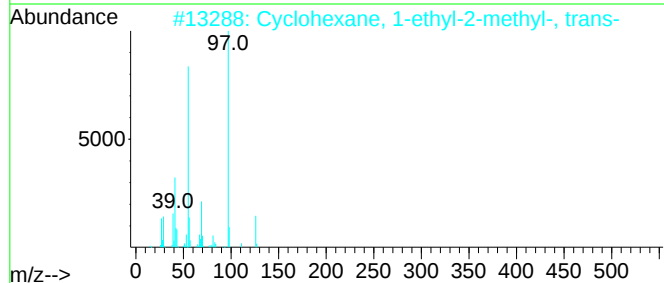
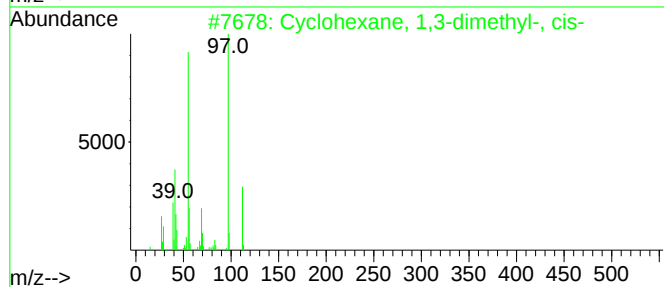
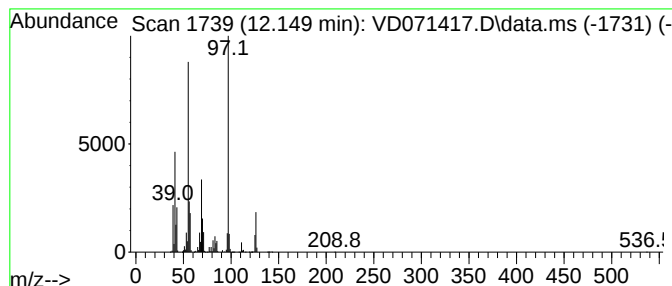
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, 1,3-dimethyl-,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.149	496.74 ug/l	17362200	Chlorobenzene-d5	11.638	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	72
2	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	72
3	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	72
4	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	72
5	2-Hexene, 3,4,4-trimethyl-	126	C9H18	053941-19-8	68



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BLIND-DUPLICATE

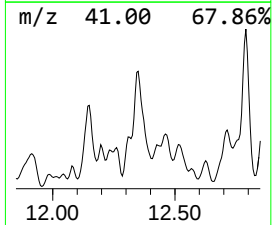
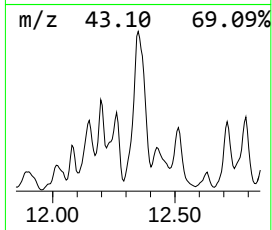
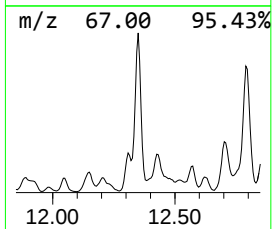
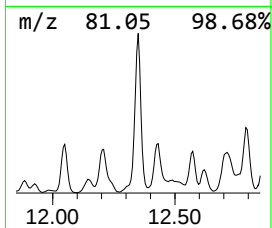
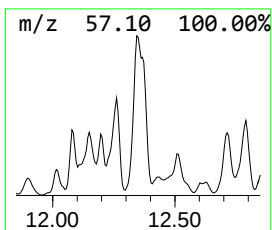
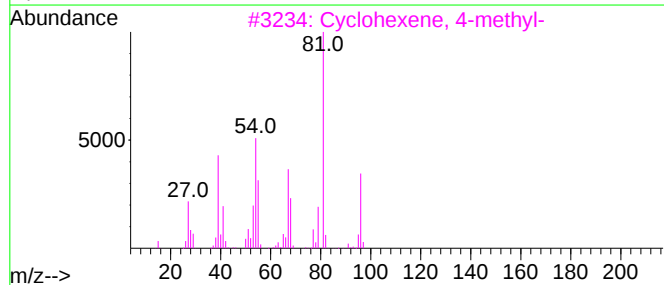
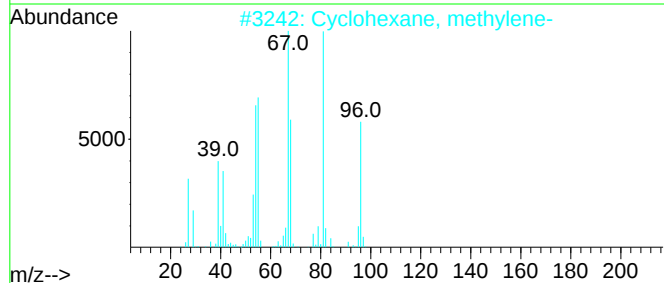
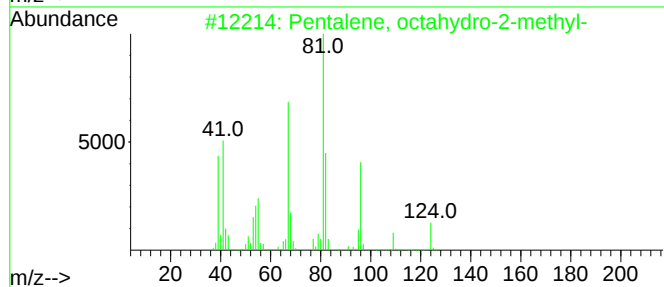
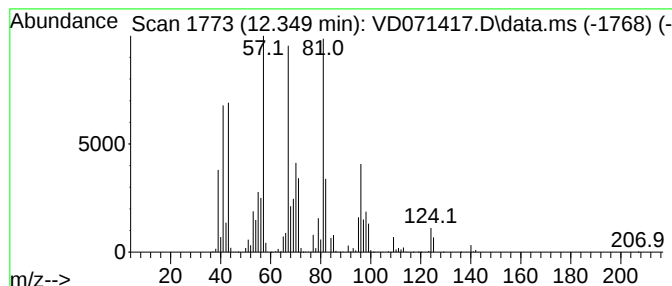
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Quant Title : SW846 8260

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

Peak Number 4 Pentalene, octahydro-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.349	600.03 ug/l	20972300	Chlorobenzene-d5	11.638

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentalene, octahydro-2-methyl-	124	C9H16	003868-64-2	60
2			Cyclohexane, methylene-	96	C7H12	001192-37-6	52
3			Cyclohexene, 4-methyl-	96	C7H12	000591-47-9	50
4			Bicyclo[3.3.1]nonane	124	C9H16	000280-65-9	47
5			Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD122121\
Data File : VD071417.D
Acq On : 21 Dec 2021 18:16
Operator : VA/SY
Sample : M5182-02
Misc : 7.69G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BLIND-DUPLICATE

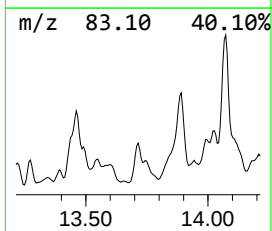
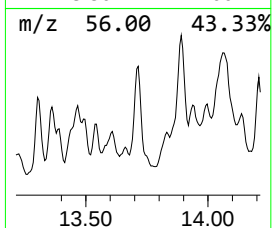
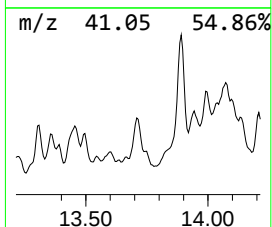
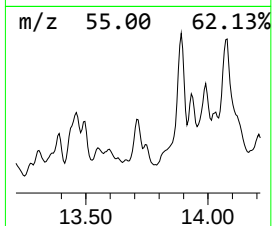
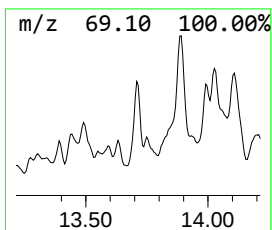
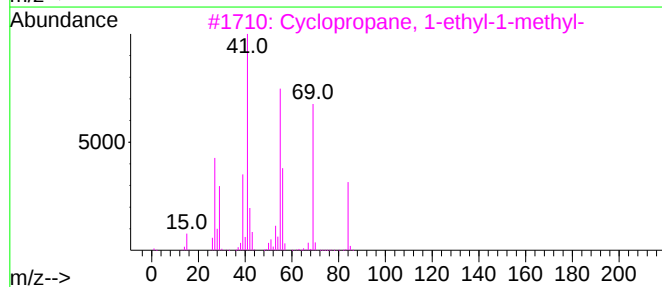
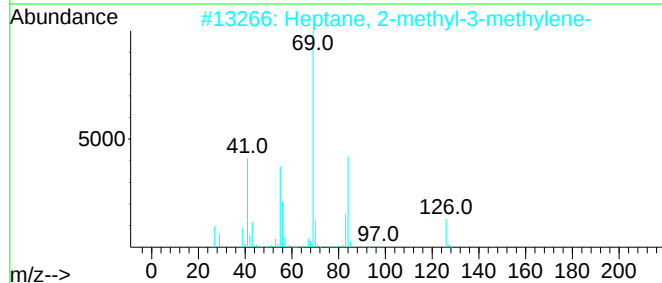
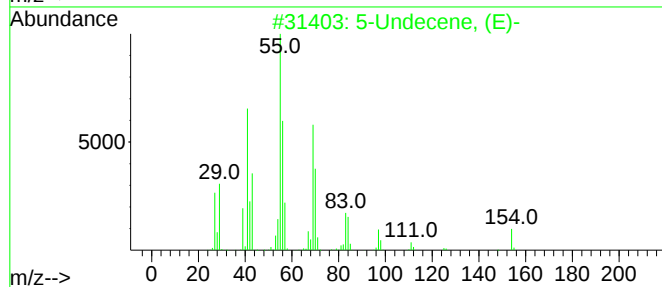
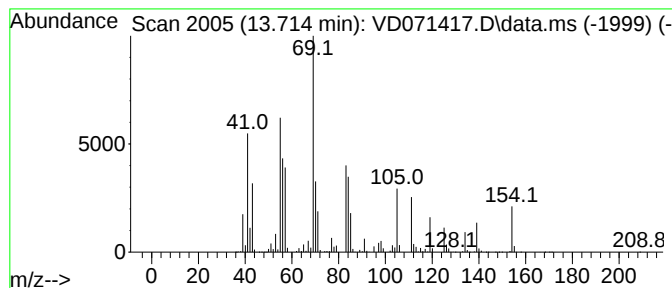
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Quant Title : SW846 8260

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

Peak Number 5 unknown13.714 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.714	167.52 ug/l	39383800	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Undecene, (E)-	154	C11H22	000764-97-6	47
2			Heptane, 2-methyl-3-methylene-	126	C9H18	062187-11-5	43
3			Cyclopropane, 1-ethyl-1-methyl-	84	C6H12	053778-43-1	38
4			Pentane, 3-methylene-	84	C6H12	000760-21-4	35
5			2-Pentene, 4-methyl-	84	C6H12	004461-48-7	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD122121\
Data File : VD071417.D
Acq On : 21 Dec 2021 18:16
Operator : VA/SY
Sample : M5182-02
Misc : 7.69G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BLIND-DUPLICATE

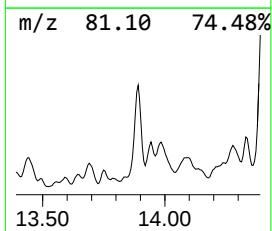
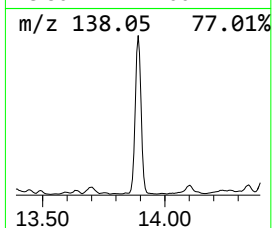
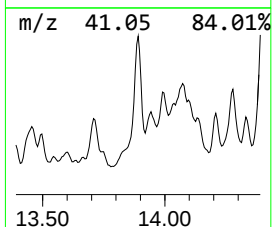
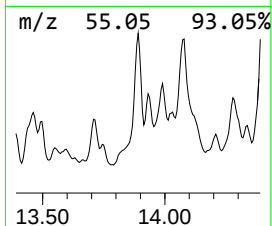
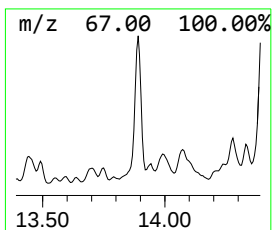
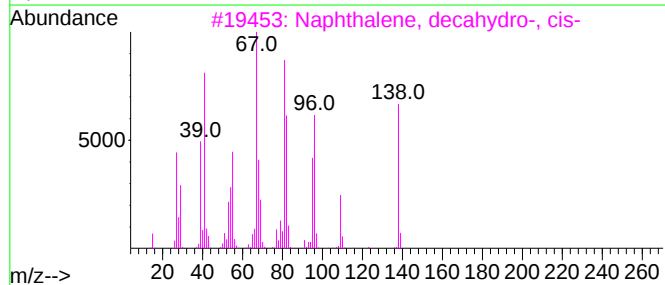
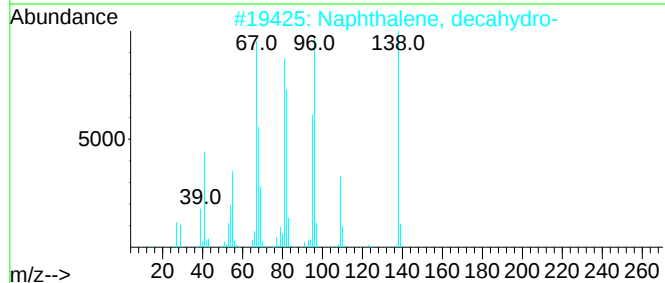
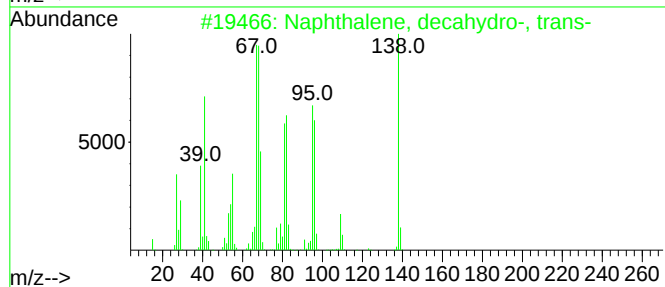
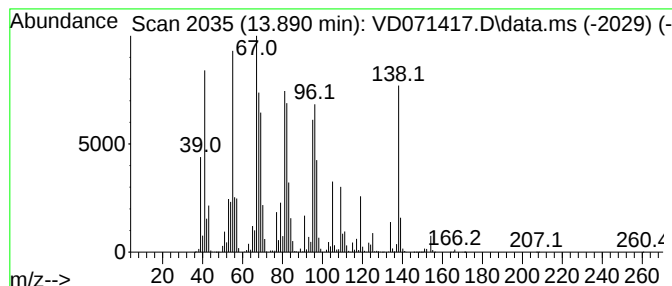
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Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, decahydro-, tr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.890	313.69 ug/l	73749800	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	96
2			Naphthalene, decahydro-	138	C10H18	000091-17-8	96
3			Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	95
4			Spiro[4.5]decane	138	C10H18	000176-63-6	93
5			Cyclohexane, 1-methyl-4-(1-methy...	138	C10H18	001124-25-0	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD122121\
Data File : VD071417.D
Acq On : 21 Dec 2021 18:16
Operator : VA/SY
Sample : M5182-02
Misc : 7.69G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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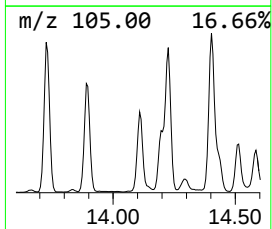
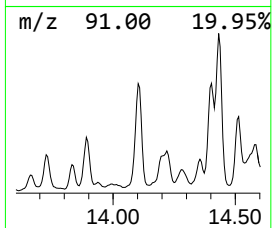
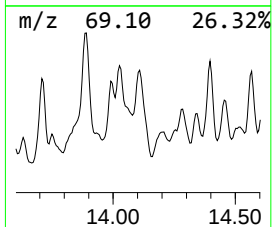
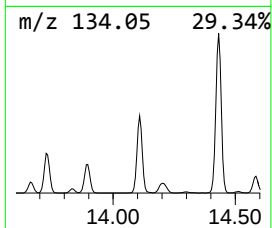
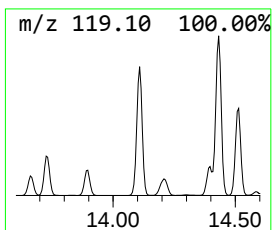
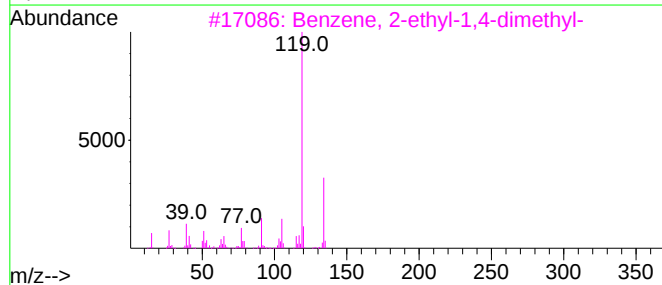
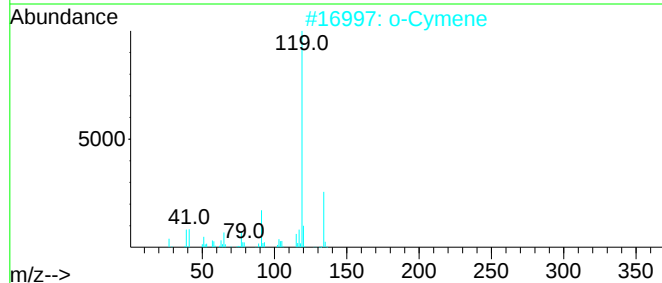
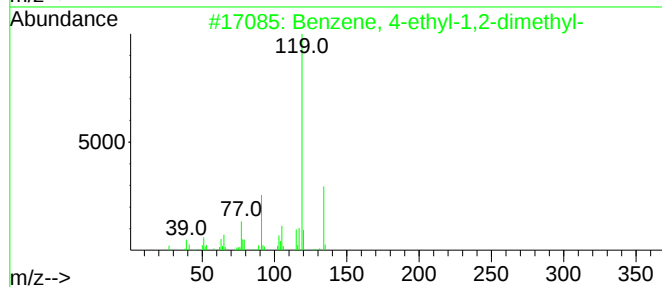
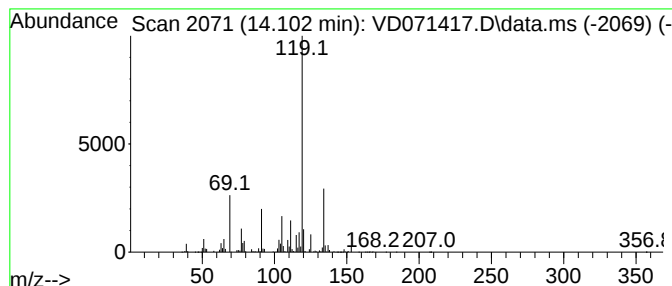
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.102	202.56 ug/l	47622500	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5			p-Cymene	134	C10H14	000099-87-6	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD122121\
Data File : VD071417.D
Acq On : 21 Dec 2021 18:16
Operator : VA/SY
Sample : M5182-02
Misc : 7.69G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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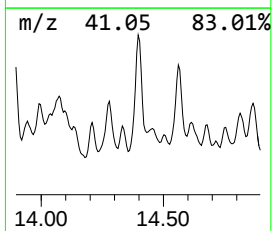
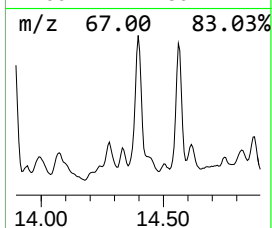
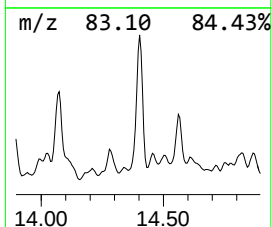
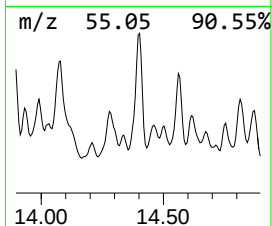
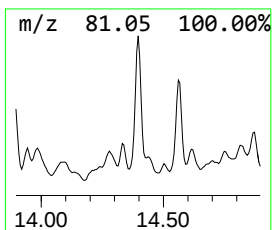
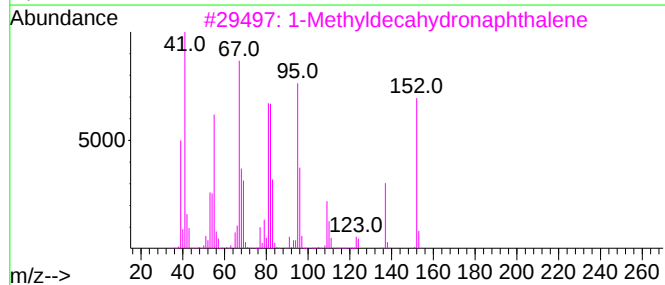
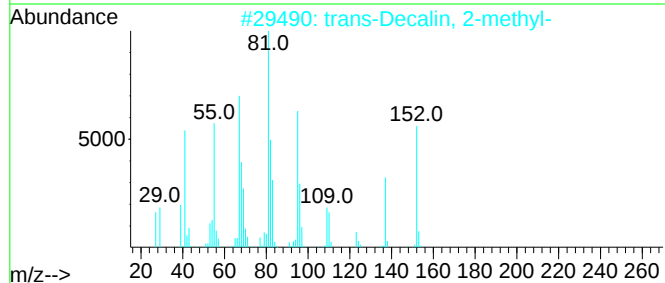
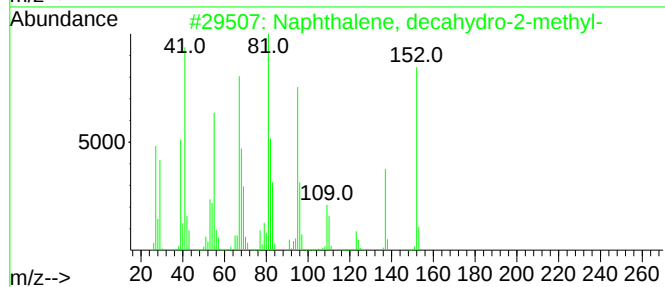
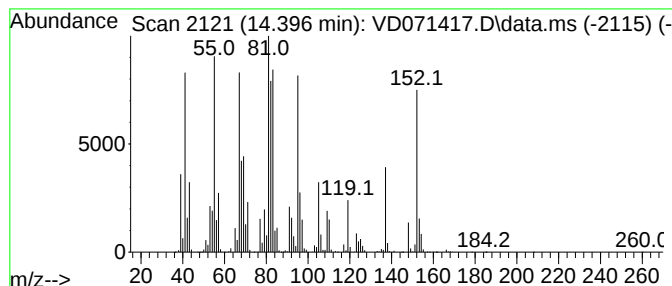
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, decahydro-2-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.396	373.47 ug/l	87805200	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	96
2			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	90
3			1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	64
4			Decalin, anti-1-methyl-, cis-	152	C11H20	1000158-89-0	52
5			Pulegone	152	C10H16O	000089-82-7	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD122121\
Data File : VD071417.D
Acq On : 21 Dec 2021 18:16
Operator : VA/SY
Sample : M5182-02
Misc : 7.69G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BLIND-DUPLICATE

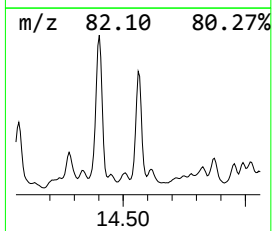
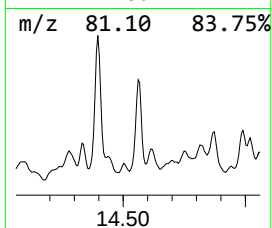
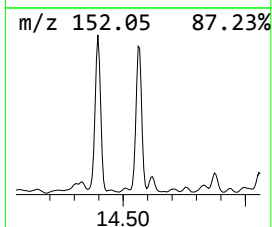
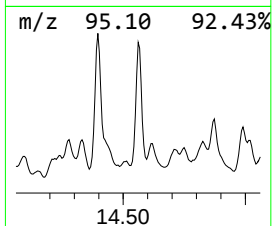
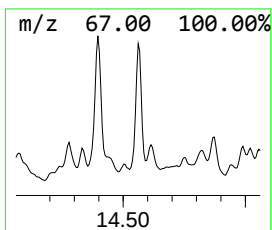
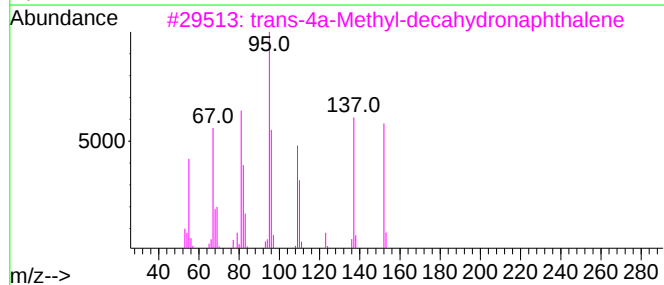
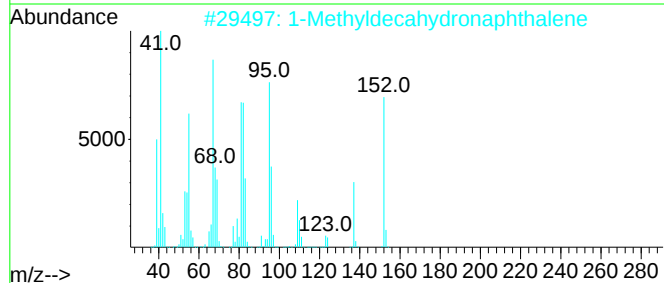
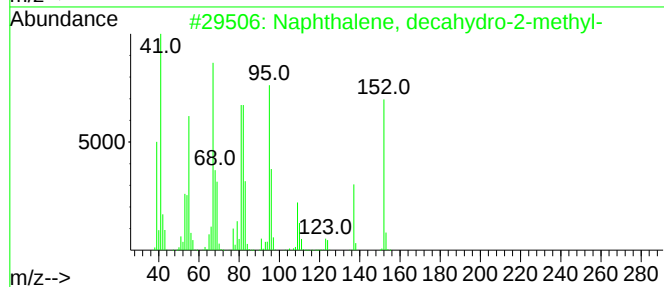
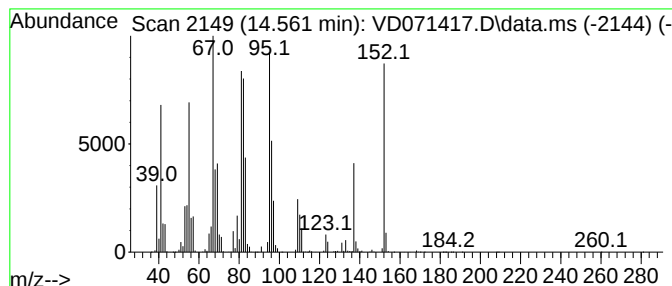
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Quant Title : SW846 8260

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

Peak Number 9 1-Methyldecahydronaphthalene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.561	250.19 ug/l	58821000	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	98
2			1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	98
3			trans-4a-Methyl-decahydronaphtha...	152	C11H20	002547-27-5	83
4			Bicyclo[3.3.2]decan-9-one	152	C10H16O	028054-91-3	64
5			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD122121\
Data File : VD071417.D
Acq On : 21 Dec 2021 18:16
Operator : VA/SY
Sample : M5182-02
Misc : 7.69G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BLIND-DUPLICATE

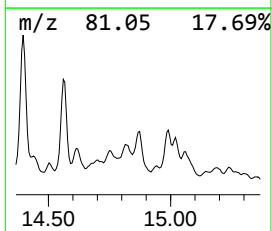
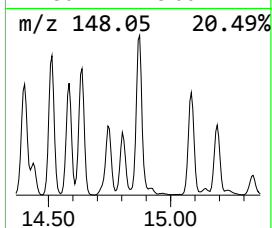
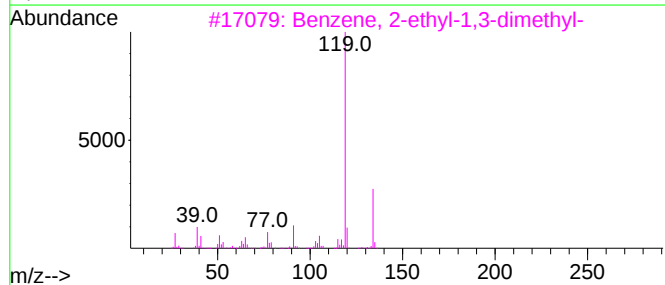
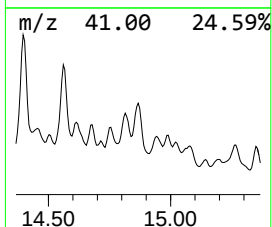
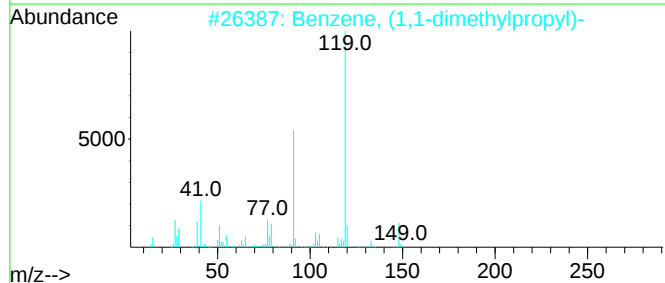
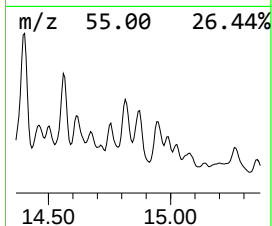
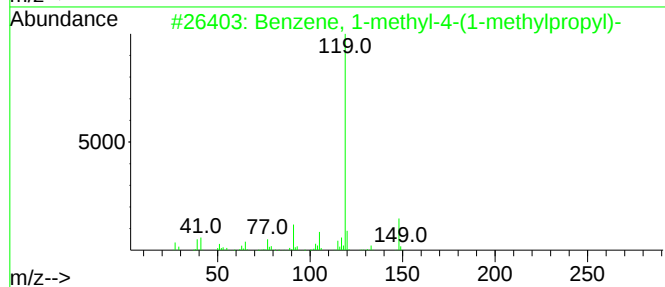
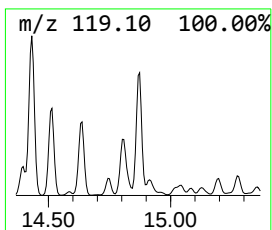
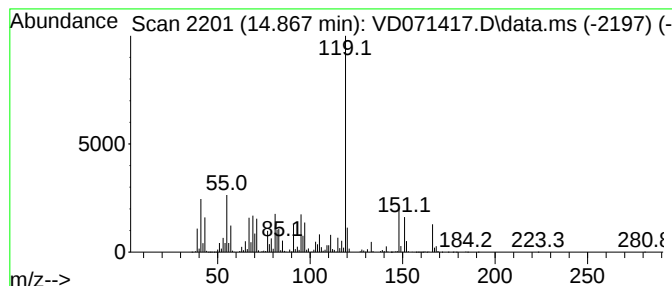
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-methyl-4-(1-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.867	198.38 ug/l	46640700	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	92
2			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	62
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	55
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	55
5			Acetic acid, (2,4-xylyl)-	164	C10H12O2	006331-04-0	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD122121\
Data File : VD071417.D
Acq On : 21 Dec 2021 18:16
Operator : VA/SY
Sample : M5182-02
Misc : 7.69G/5.00ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BLIND-DUPLICATE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclohexane, 1,...	11.444	187.6	ug/l	6558380	3	11.638	1747600	50.0
Cyclohexane, 1-...	11.920	330.9	ug/l	11566700	3	11.638	1747600	50.0
Cyclohexane, 1,...	12.149	496.7	ug/l	17362200	3	11.638	1747600	50.0
Pentalene, octa...	12.349	600.0	ug/l	20972300	3	11.638	1747600	50.0
unknown13.714	13.714	167.5	ug/l	39383800	4	13.567	11755200	50.0
Naphthalene, de...	13.890	313.7	ug/l	73749800	4	13.567	11755200	50.0
Benzene, 4-ethy...	14.102	202.6	ug/l	47622500	4	13.567	11755200	50.0
Naphthalene, de...	14.396	373.5	ug/l	87805200	4	13.567	11755200	50.0
1-Methyldecahyd...	14.561	250.2	ug/l	58821000	4	13.567	11755200	50.0
Benzene, 1-meth...	14.867	198.4	ug/l	46640700	4	13.567	11755200	50.0