

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD050721\
 Data File : VD069113.D
 Acq On : 07 May 2021 20:18
 Operator : VA/SY
 Sample : M2295-01
 Misc : 3.88G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 DRUM-A-B-C

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\82D050421S.M
 Title : SW846 8260

Signal : TIC: VD069113.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.226	45	52	55	rBV3	5192	12827	1.24%	0.104%
2	2.491	93	97	101	rBV	10202	14963	1.45%	0.122%
3	3.573	272	281	295	rBV	100353	258833	25.01%	2.106%
4	4.156	371	380	392	rBV2	37681	104452	10.09%	0.850%
5	4.444	419	429	430	rBV4	5730	12734	1.23%	0.104%
6	4.962	507	517	526	rBV3	16739	55550	5.37%	0.452%
7	7.914	1010	1019	1024	rBV2	67523	165175	15.96%	1.344%
8	7.979	1024	1030	1039	rVB2	172314	389034	37.58%	3.166%
9	8.332	1077	1090	1100	rBV	122603	274389	26.51%	2.233%
10	8.861	1172	1180	1189	rBV	275018	550513	53.19%	4.480%
11	9.426	1271	1276	1283	rVB5	7066	14413	1.39%	0.117%
12	9.973	1363	1369	1373	rBV5	6249	11992	1.16%	0.098%
13	10.114	1384	1393	1397	rBV4	7790	15401	1.49%	0.125%
14	10.338	1421	1431	1445	rBV	494760	912729	88.18%	7.427%
15	10.520	1455	1462	1467	rVB7	15324	34696	3.35%	0.282%
16	10.679	1484	1489	1494	rVB5	15697	23951	2.31%	0.195%
17	10.773	1498	1505	1510	rBV6	21654	41149	3.98%	0.335%
18	10.867	1515	1521	1526	rVB3	10993	20141	1.95%	0.164%
19	10.950	1526	1535	1541	rBV4	10890	20528	1.98%	0.167%
20	11.085	1554	1558	1561	rBV4	12804	15216	1.47%	0.124%
21	11.120	1561	1564	1568	rVV4	21647	34575	3.34%	0.281%
22	11.191	1572	1576	1580	rVV2	30207	46306	4.47%	0.377%
23	11.244	1580	1585	1590	rVB3	28144	55938	5.40%	0.455%
24	11.297	1590	1594	1598	rVB6	13640	20750	2.00%	0.169%
25	11.350	1598	1603	1608	rBV5	19351	33167	3.20%	0.270%
26	11.444	1608	1619	1628	rVB6	86803	257515	24.88%	2.095%
27	11.526	1628	1633	1638	rBV	45723	71014	6.86%	0.578%
28	11.644	1646	1653	1664	rVB	460380	805237	77.79%	6.552%
29	11.779	1665	1676	1679	rBV4	58558	126244	12.20%	1.027%
30	11.826	1679	1684	1687	rVV4	98340	180763	17.46%	1.471%
31	11.879	1687	1693	1698	rVV3	144383	330431	31.92%	2.689%
32	11.926	1698	1701	1707	rVB	101171	164923	15.93%	1.342%
33	11.985	1707	1711	1716	rBV7	9964	18258	1.76%	0.149%
34	12.067	1716	1725	1731	rVB10	13594	38268	3.70%	0.311%
35	12.144	1731	1738	1751	rBV3	156969	504004	48.69%	4.101%
36	12.261	1753	1758	1762	rVB2	41994	75594	7.30%	0.615%

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Integrator: RTE

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D050421S.M
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37	12.314	1762	1767	1769	rBV3	27217	46849	4.53%	0.381%
38	12.349	1769	1773	1776	rBV5	63009	107134	10.35%	0.872%
39	12.397	1776	1781	1785	rVV4	73094	120852	11.68%	0.983%
40	12.555	1805	1808	1810	rBV3	14906	17929	1.73%	0.146%
41	12.626	1814	1820	1829	rBV	625062	1035090	100.00%	8.423%
42	12.714	1829	1835	1839	rBV8	35848	54518	5.27%	0.444%
43	12.797	1844	1849	1855	rVB5	92189	169261	16.35%	1.377%
44	12.873	1855	1862	1865	rBV6	45813	99601	9.62%	0.810%
45	12.949	1865	1875	1881	rBV4	165935	374193	36.15%	3.045%
46	13.002	1881	1884	1891	rVB5	56970	83429	8.06%	0.679%
47	13.085	1895	1898	1903	rVB5	49528	83541	8.07%	0.680%
48	13.185	1903	1915	1924	rBV6	126840	375854	36.31%	3.058%
49	13.273	1925	1930	1933	rBV5	29618	54422	5.26%	0.443%
50	13.396	1947	1951	1954	rBV4	42492	62115	6.00%	0.505%
51	13.455	1955	1961	1966	rVV5	106413	214801	20.75%	1.748%
52	13.567	1972	1980	1989	rVB	508430	923990	89.27%	7.519%
53	13.661	1989	1996	1999	rBV2	74346	156711	15.14%	1.275%
54	13.761	2011	2013	2020	rVB5	53893	87936	8.50%	0.716%
55	13.891	2030	2035	2041	rBV3	208343	401903	38.83%	3.270%
56	13.991	2048	2052	2056	rBV5	40041	73774	7.13%	0.600%
57	14.102	2067	2071	2077	rVB6	41924	84132	8.13%	0.685%
58	14.226	2086	2092	2096	rBV7	50751	107897	10.42%	0.878%
59	14.273	2096	2100	2106	rBV5	48248	87066	8.41%	0.708%
60	14.338	2107	2111	2116	rBV5	42297	72651	7.02%	0.591%
61	14.402	2116	2122	2126	rBV5	123086	257746	24.90%	2.097%
62	14.567	2138	2150	2153	rBV4	71173	188197	18.18%	1.531%
63	14.620	2155	2159	2165	rVB8	42816	78635	7.60%	0.640%
64	14.814	2187	2192	2198	rBV6	88638	203963	19.70%	1.660%
65	14.867	2198	2201	2207	rVB8	63641	123267	11.91%	1.003%
66	15.020	2225	2227	2233	rVB5	39965	58252	5.63%	0.474%
67	15.085	2235	2238	2243	rVB6	64494	106980	10.34%	0.871%
68	15.202	2253	2258	2263	rVB5	41529	77602	7.50%	0.631%
69	15.390	2286	2290	2295	rVB	65858	103899	10.04%	0.845%
70	15.667	2332	2337	2338	rBV5	22976	35079	3.39%	0.285%
71	15.849	2364	2368	2372	rBV7	17848	37038	3.58%	0.301%
72	16.208	2423	2429	2435	rBV7	35145	76200	7.36%	0.620%
73	16.479	2470	2475	2482	rBV3	49022	112435	10.86%	0.915%
74	16.685	2504	2510	2520	rVB3	89560	222855	21.53%	1.813%

Sum of corrected areas: 12289470

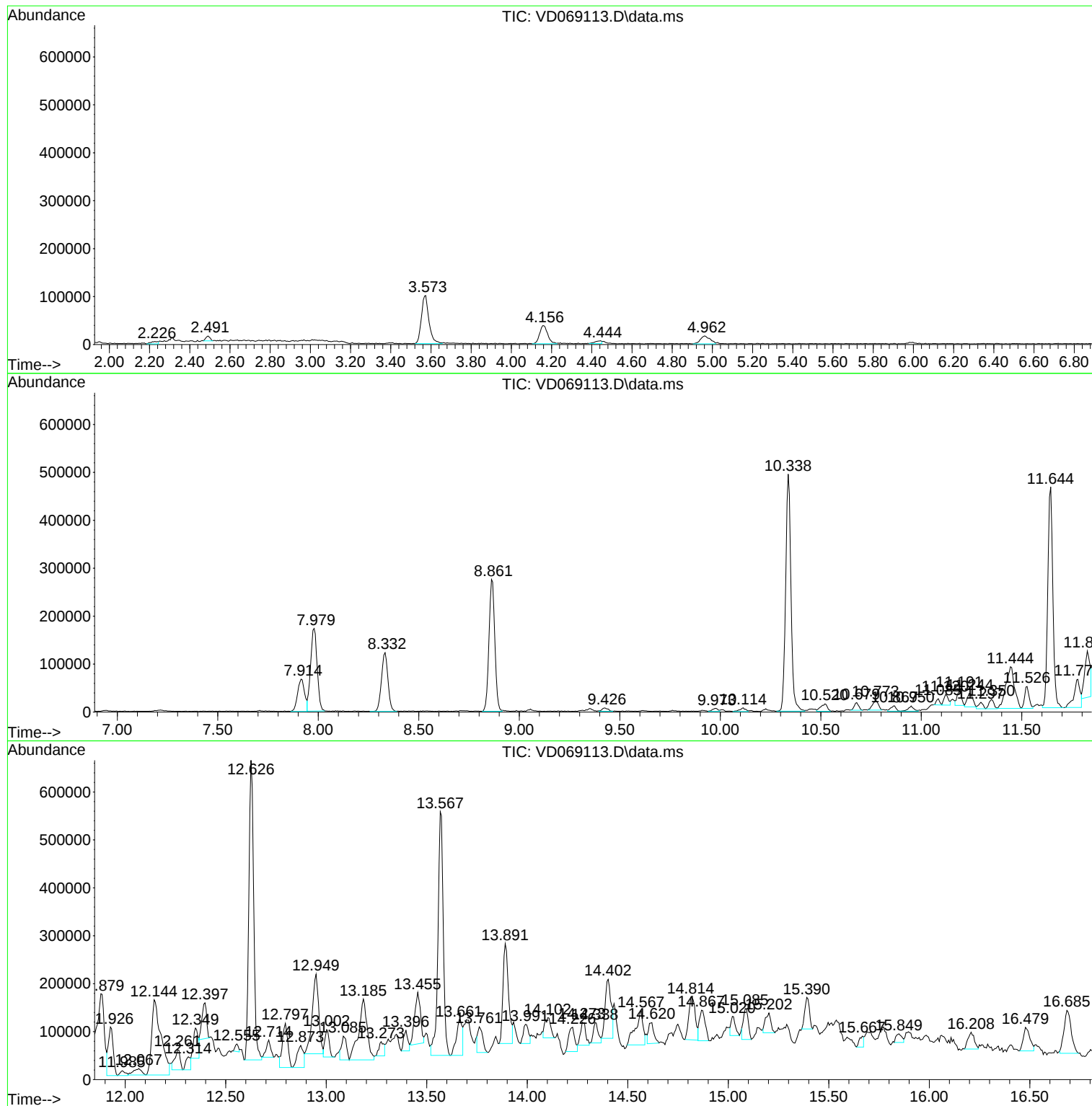
Instrument :
MSVOA_D
ClientSampleId :
DRUM-A-B-C

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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D050421S.M
Quant Title : SW846 8260

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



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Operator : VA/SY
Sample : M2295-01
Misc : 3.88G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DRUM-A-B-C

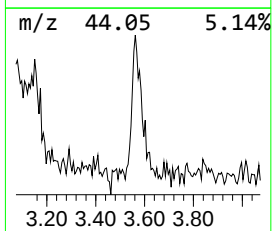
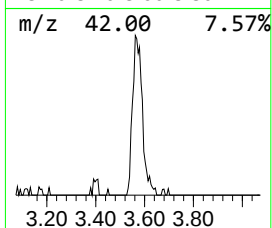
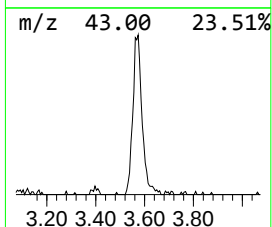
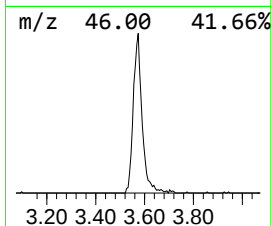
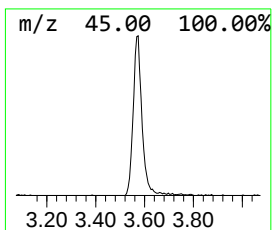
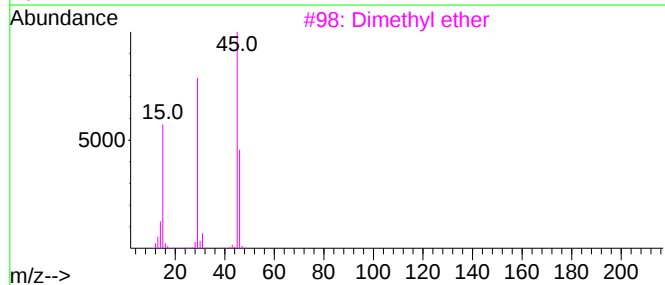
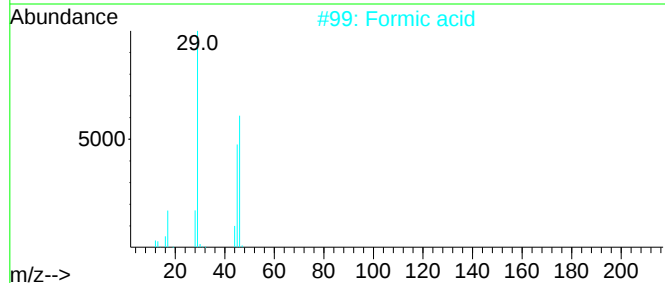
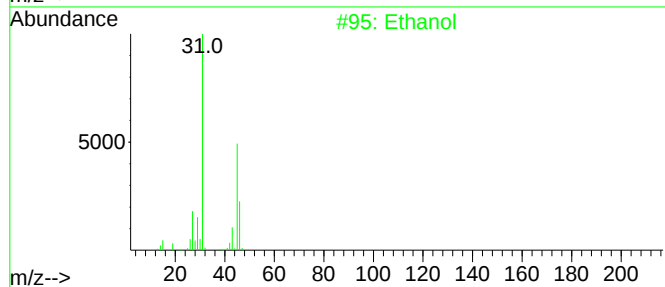
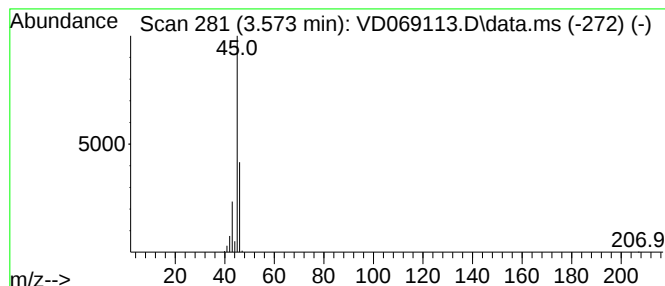
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D050421S.M
Quant Title : SW846 8260

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

Peak Number 1 Ethanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.573	33.27 ug/l	258833	Pentafluorobenzene	7.973

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol			46	C2H6O	000064-17-5	64
2	Formic acid			46	CH2O2	000064-18-6	5
3	Dimethyl ether			46	C2H6O	000115-10-6	5
4	Oxalic acid			90	C2H2O4	000144-62-7	4
5	Methane, nitroso-			45	CH3NO	000865-40-7	3



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ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DRUM-A-B-C

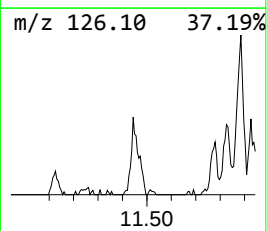
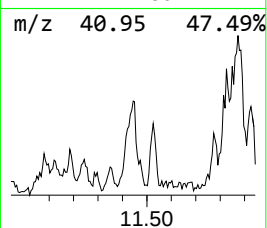
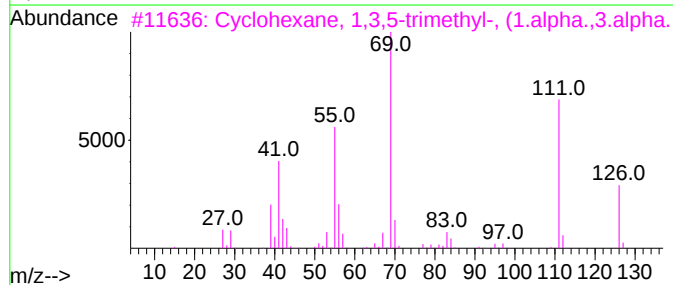
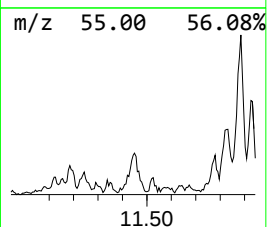
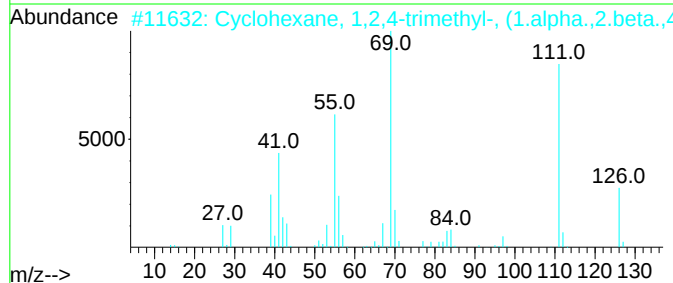
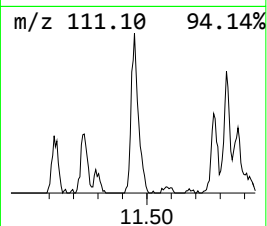
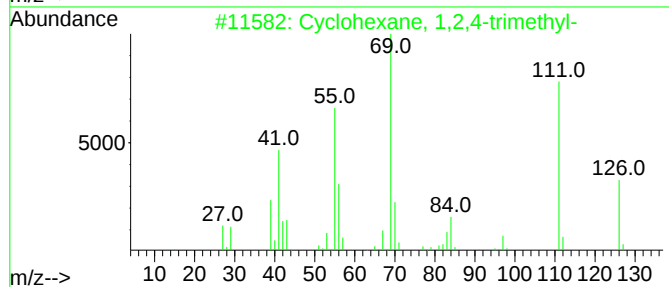
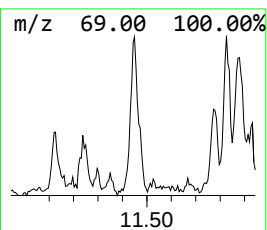
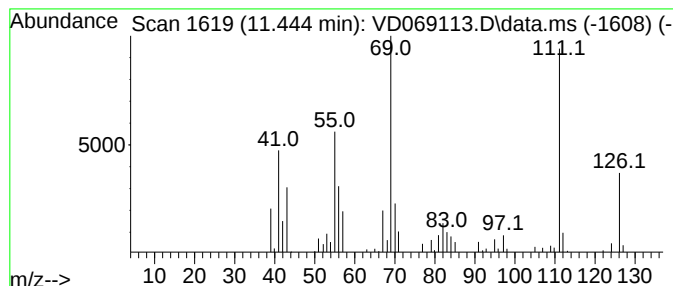
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D050421S.M
Quant Title : SW846 8260

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

Peak Number 2 Cyclohexane, 1,2,4-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.444	15.99 ug/l	257515	Chlorobenzene-d5	11.644

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	93
2			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	90
3			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	83
4			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	72
5			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	72



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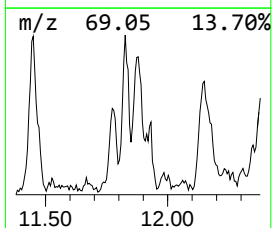
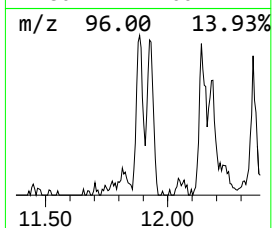
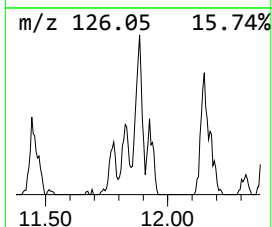
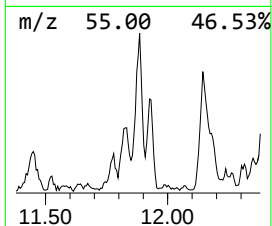
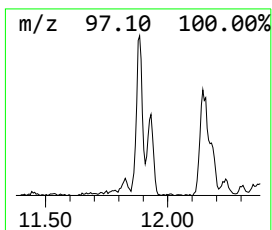
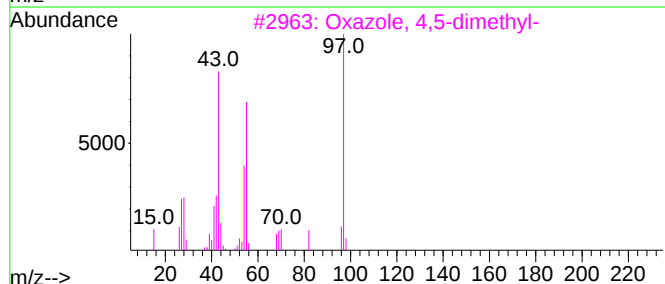
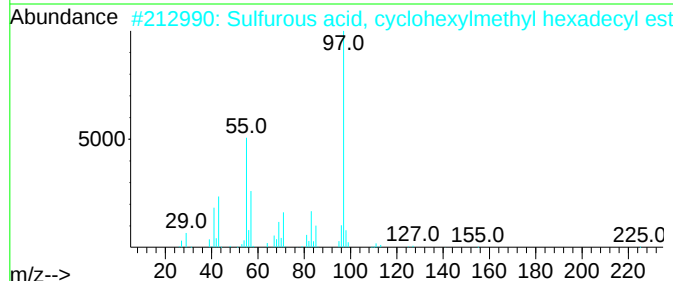
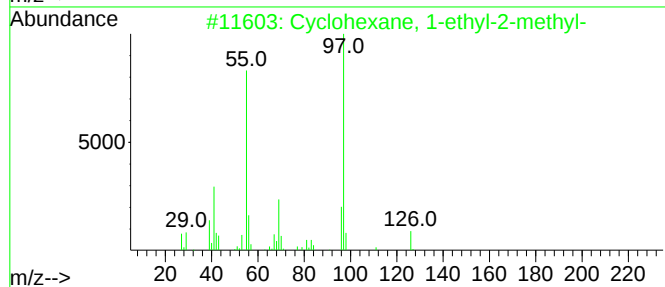
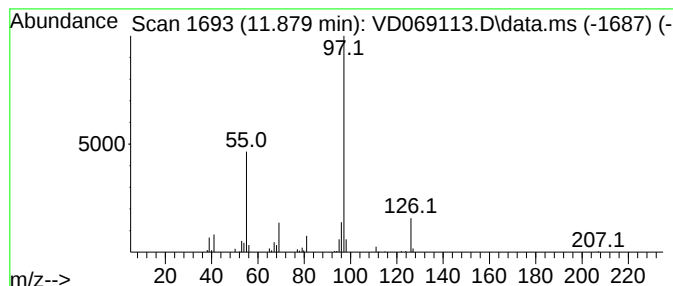
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.879	20.52 ug/l	330431	Chlorobenzene-d5	11.644	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	87
2	Sulfurous acid, cyclohexylmethyl...	402	C23H46O3S	1000309-22-4	78
3	Oxazole, 4,5-dimethyl-	97	C5H7NO	020662-83-3	64
4	Furan, 2,3-dihydro-4-(1-methylpr...	126	C8H14O	034379-54-9	64
5	1H-Pyrazole, 4,5-dihydro-3-methy...	126	C7H14N2	026964-49-8	64



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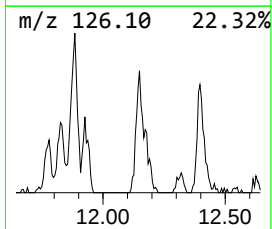
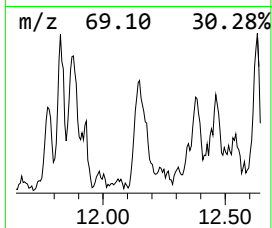
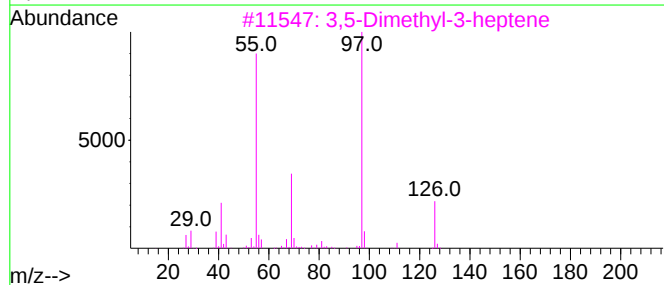
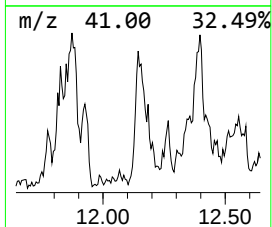
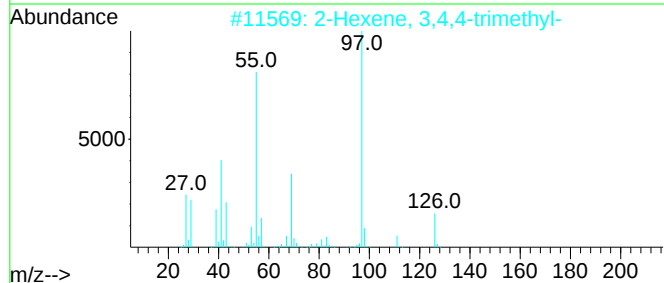
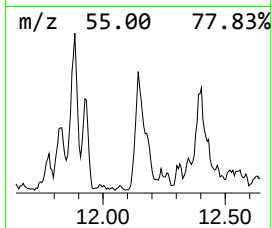
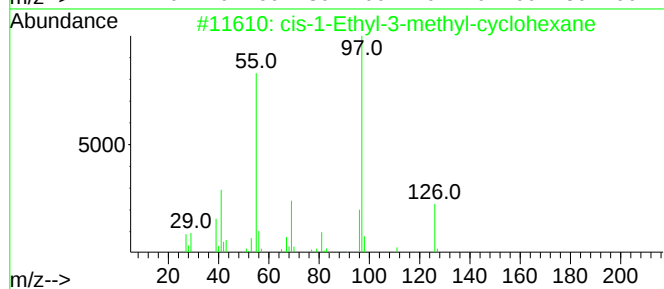
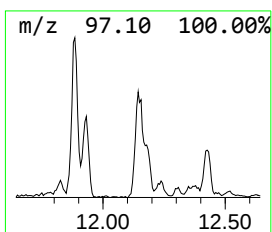
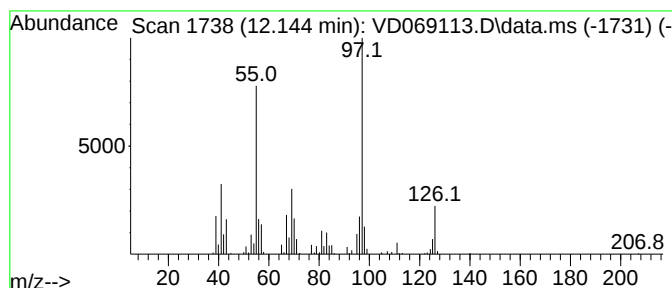
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Quant Title : SW846 8260

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

Peak Number 4 cis-1-Ethyl-3-methyl-cyclohexane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.144	31.30 ug/l	504004	Chlorobenzene-d5	11.644

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	81
2		2-Hexene, 3,4,4-trimethyl-	126	C9H18	053941-19-8	74
3		3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	72
4		1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	72
5		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD050721\
Data File : VD069113.D
Acq On : 07 May 2021 20:18
Operator : VA/SY
Sample : M2295-01
Misc : 3.88G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DRUM-A-B-C

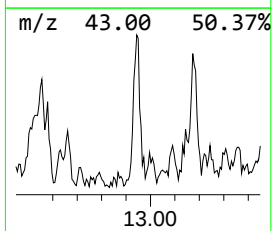
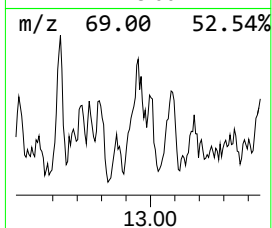
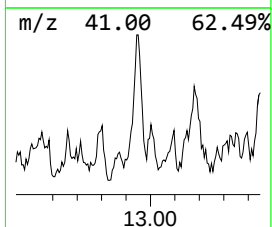
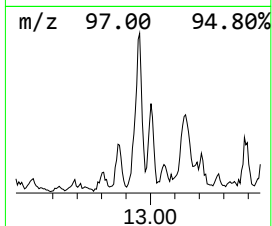
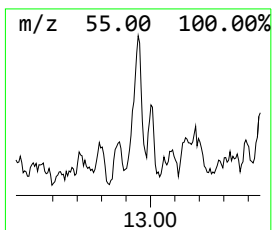
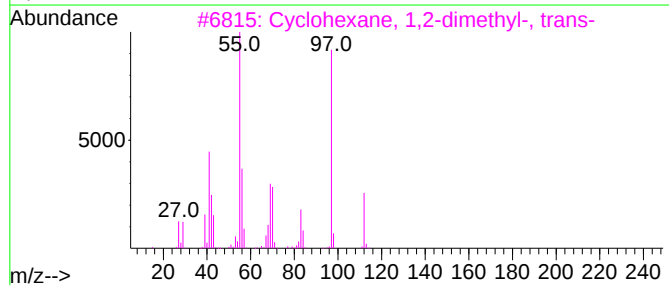
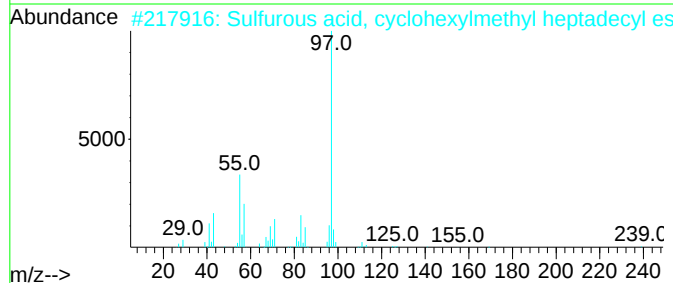
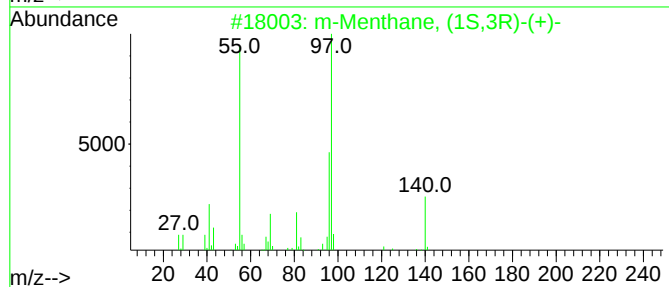
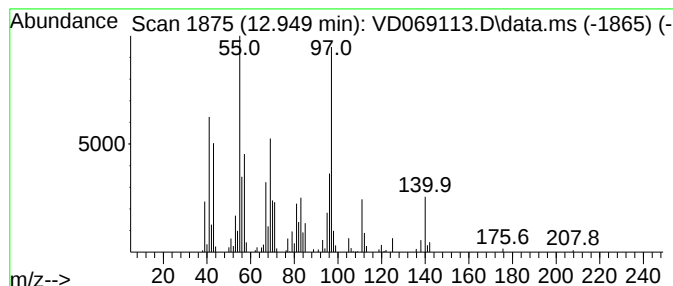
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Quant Title : SW846 8260

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

Peak Number 5 m-Menthane, (1S,3R)-(+)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.950	20.25 ug/l	374193	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			m-Menthane, (1S,3R)-(+)-	140	C10H20	013837-66-6	58
2			Sulfurous acid, cyclohexylmethyl...	416	C24H48O3S	1000309-22-5	52
3			Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	49
4			Cyclopentane, 1,2-dimethyl-3-(1-...	140	C10H20	000489-20-3	47
5			Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD050721\
Data File : VD069113.D
Acq On : 07 May 2021 20:18
Operator : VA/SY
Sample : M2295-01
Misc : 3.88G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DRUM-A-B-C

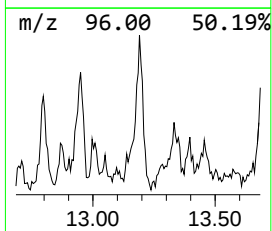
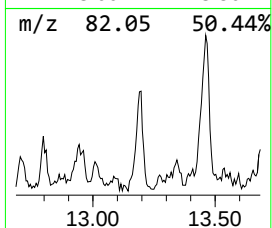
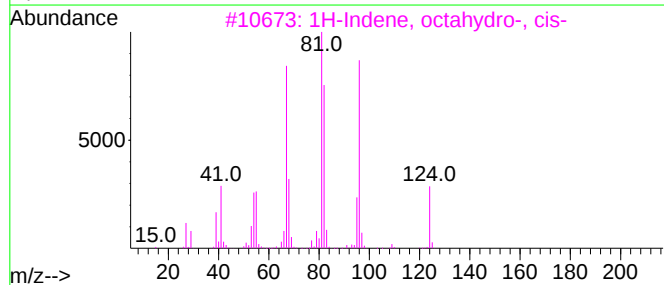
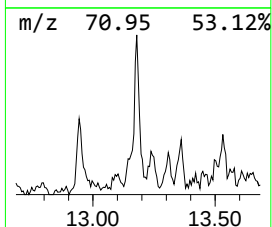
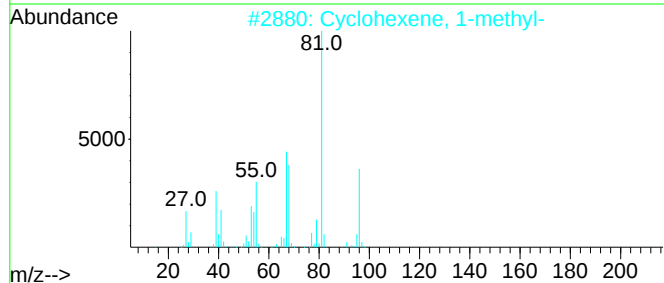
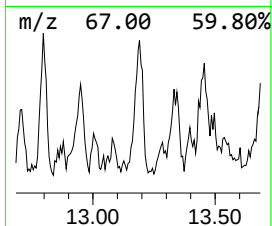
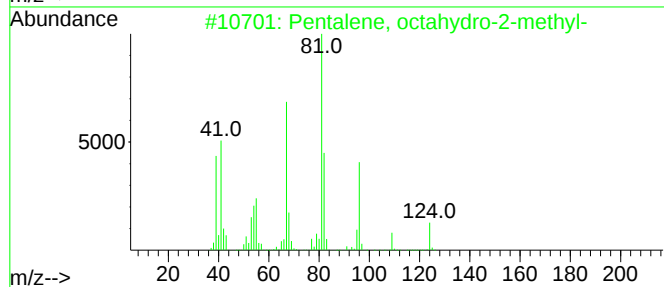
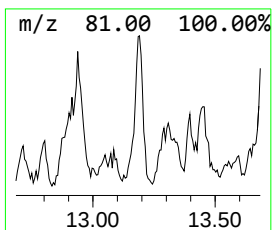
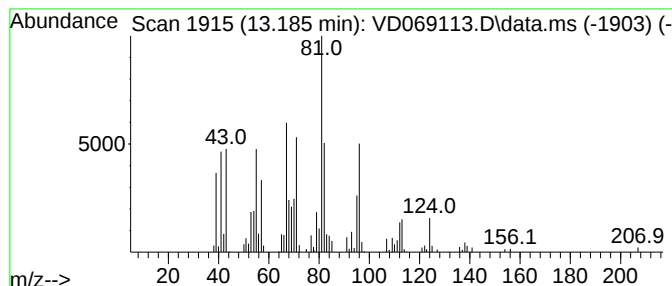
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Quant Title : SW846 8260

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

Peak Number 6 Pentalene, octahydro-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.185	20.34 ug/l	375854	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentalene, octahydro-2-methyl-	124	C9H16	003868-64-2	58
2			Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	55
3			1H-Indene, octahydro-, cis-	124	C9H16	004551-51-3	50
4			Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	50
5			Cyclohexene,3-propyl-	124	C9H16	003983-06-0	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD050721\
Data File : VD069113.D
Acq On : 07 May 2021 20:18
Operator : VA/SY
Sample : M2295-01
Misc : 3.88G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DRUM-A-B-C

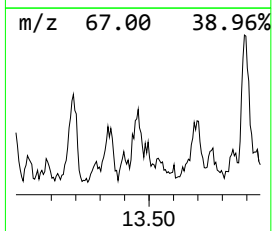
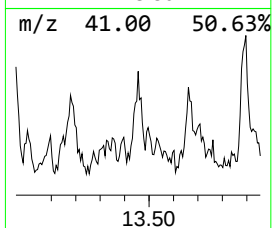
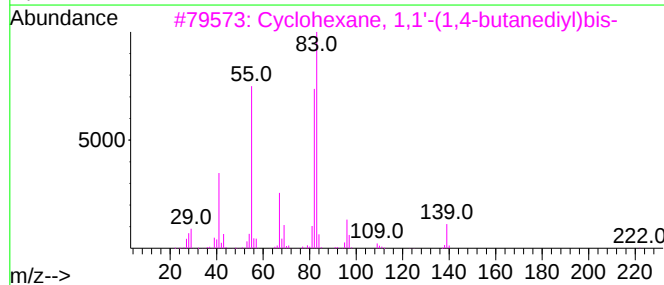
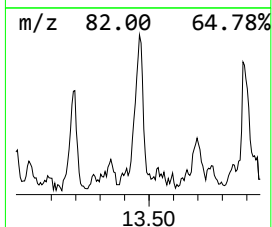
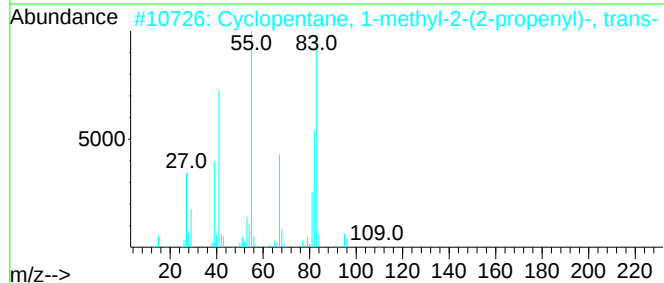
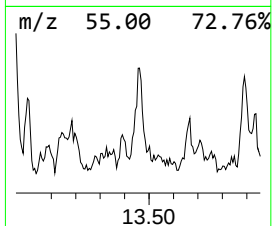
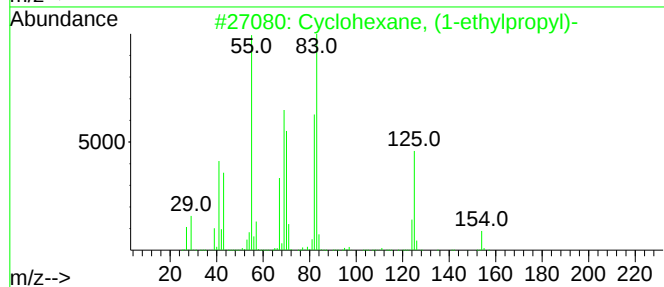
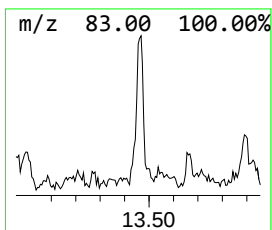
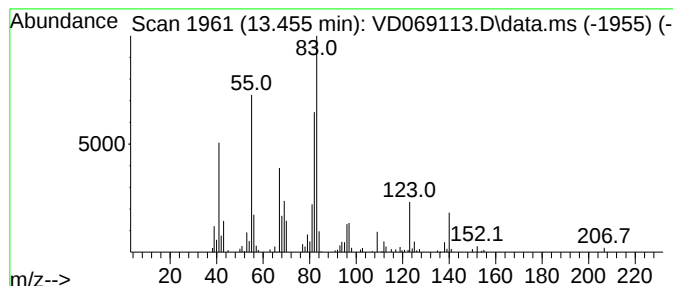
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Quant Title : SW846 8260

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

Peak Number 7 Cyclohexane, (1-ethylpropyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.455	11.62 ug/l	214801	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (1-ethylpropyl)-	154	C11H22	026321-98-2	53
2			Cyclopentane, 1-methyl-2-(2-prop...	124	C9H16	050746-53-7	53
3			Cyclohexane, 1,1'-(1,4-butanedi...	222	C16H30	006165-44-2	53
4			Cyclohexane, butyl-	140	C10H20	001678-93-9	52
5			Cyclohexane, (4-methylpentyl)-	168	C12H24	061142-20-9	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD050721\
Data File : VD069113.D
Acq On : 07 May 2021 20:18
Operator : VA/SY
Sample : M2295-01
Misc : 3.88G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DRUM-A-B-C

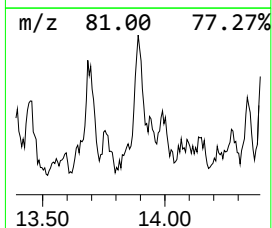
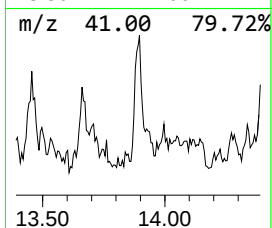
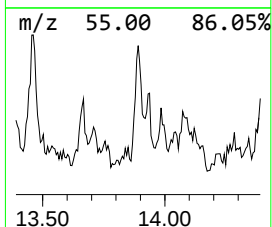
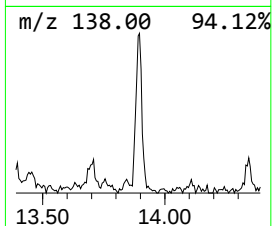
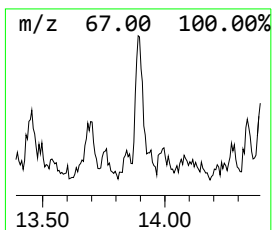
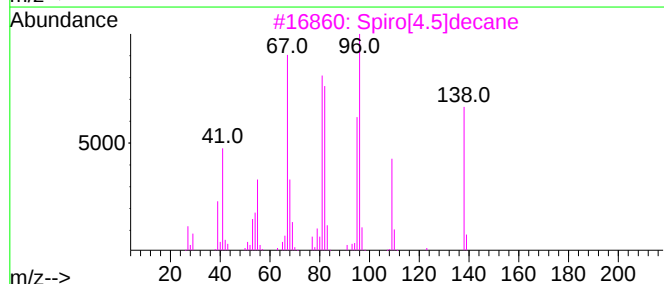
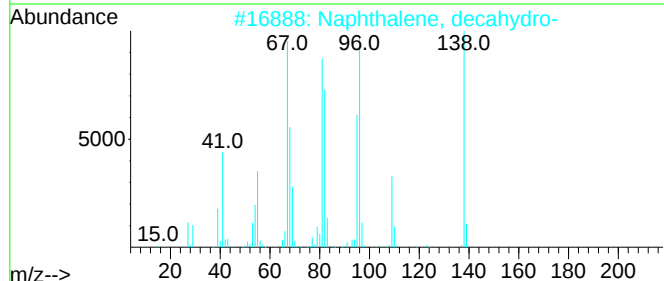
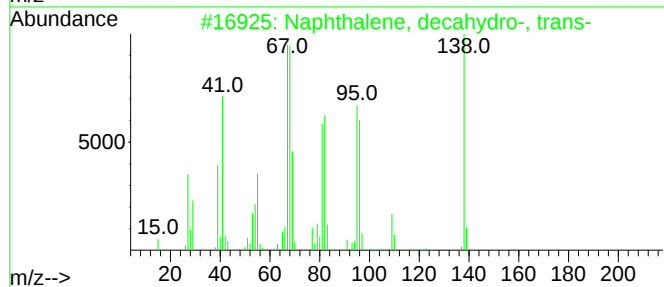
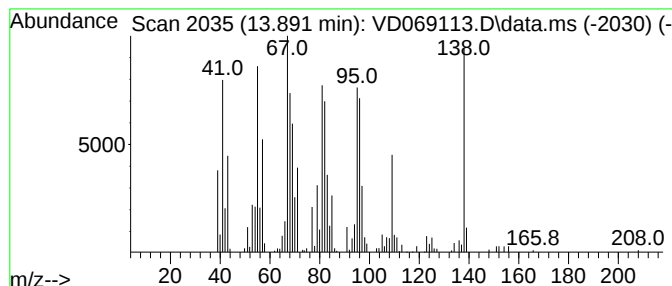
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D050421S.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, decahydro-, tr... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.891	21.75 ug/l	401903	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	94
2			Naphthalene, decahydro-	138	C10H18	000091-17-8	90
3			Spiro[4.5]decane	138	C10H18	000176-63-6	76
4			Cyclohexane, 1-methyl-4-(1-methy...	138	C10H18	001124-25-0	68
5			Cyclodecene	138	C10H18	003618-12-0	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD050721\
Data File : VD069113.D
Acq On : 07 May 2021 20:18
Operator : VA/SY
Sample : M2295-01
Misc : 3.88G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DRUM-A-B-C

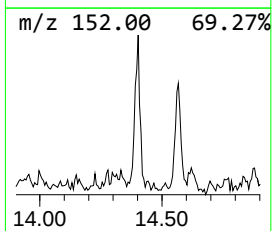
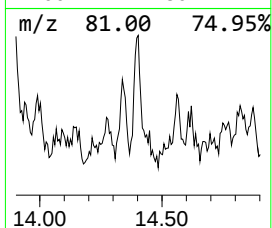
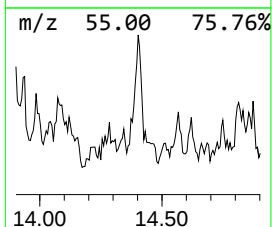
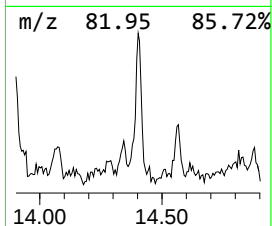
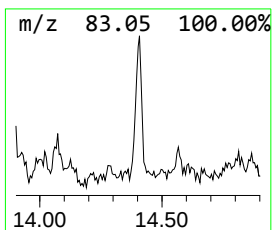
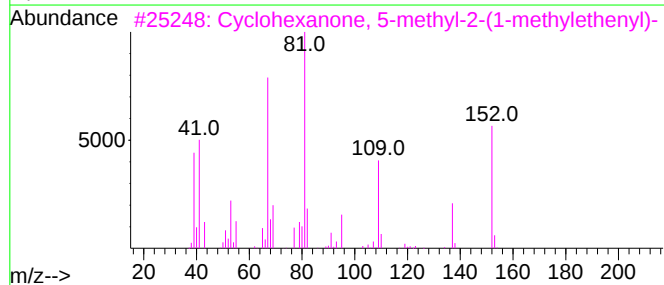
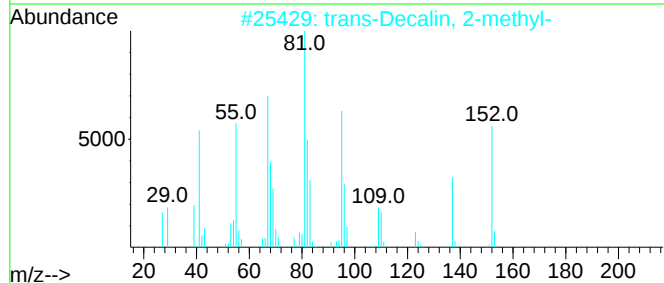
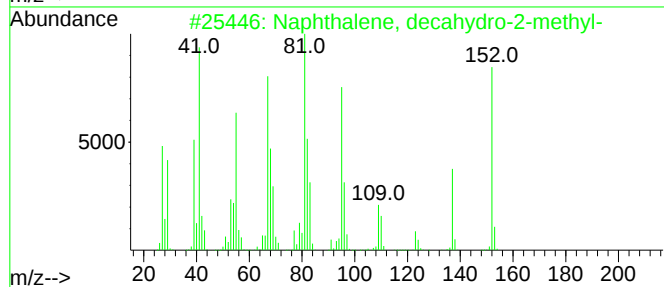
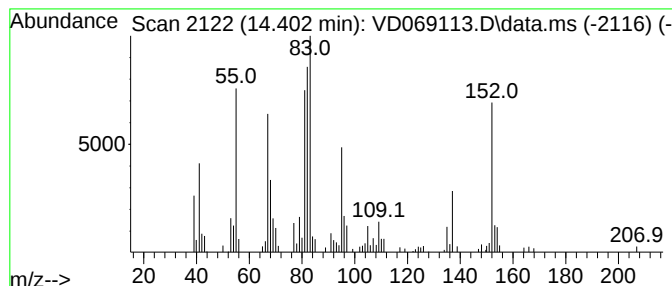
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D050421S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, decahydro-2-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.402	13.95 ug/l	257746	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	64
2			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	58
3			Cyclohexanone, 5-methyl-2-(1-met...	152	C10H16O	000529-00-0	50
4			Bicyclo[4.1.0]heptan-3-one, 4,7,...	152	C10H16O	004176-04-9	49
5			Dodeca-1,6-dien-12-ol, 6,10-dime...	210	C14H26O	1000156-13-8	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD050721\
Data File : VD069113.D
Acq On : 07 May 2021 20:18
Operator : VA/SY
Sample : M2295-01
Misc : 3.88G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DRUM-A-B-C

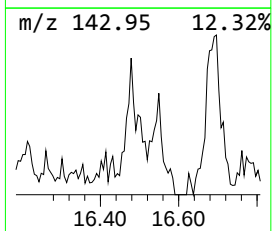
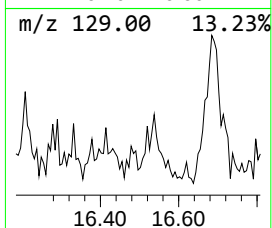
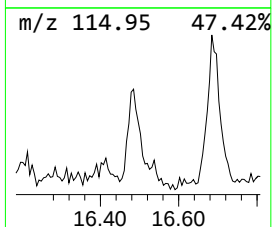
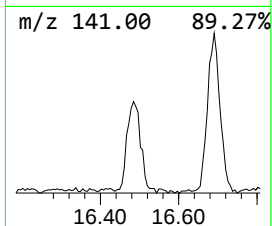
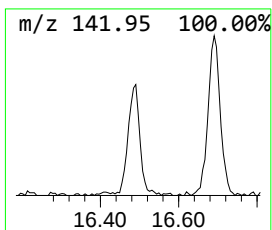
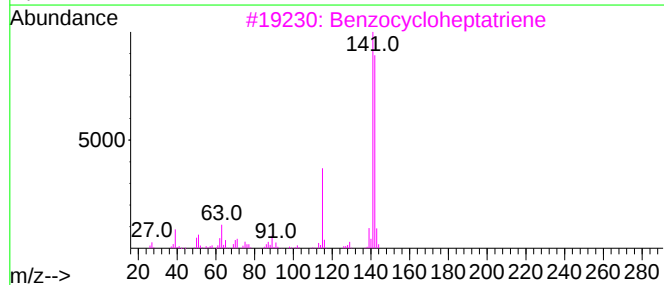
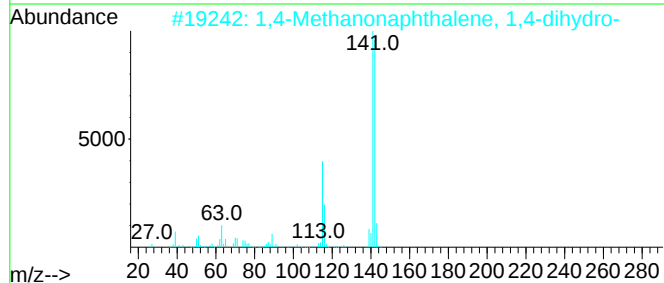
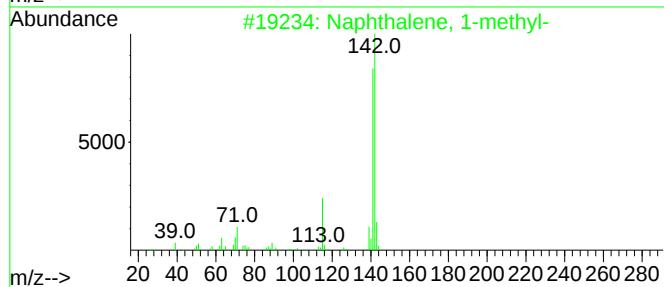
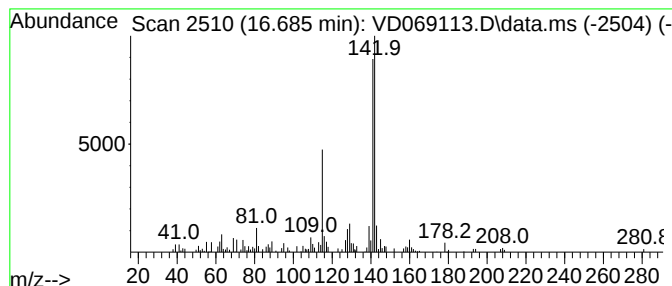
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.685	12.06 ug/l	222855	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
2			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
3			Benzocycloheptatriene	142	C11H10	000264-09-5	90
4			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	87
5			Spiro(9-methylenetricyclo[6.2.1....	184	C13H12O	033929-47-4	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD050721\
 Data File : VD069113.D
 Acq On : 07 May 2021 20:18
 Operator : VA/SY
 Sample : M2295-01
 Misc : 3.88G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 DRUM-A-B-C

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D050421S.M
 Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol	3.573	33.3	ug/l	258833	1	7.973	389034	50.0
Cyclohexane, 1,...	11.444	16.0	ug/l	257515	3	11.644	805237	50.0
Cyclohexane, 1-...	11.879	20.5	ug/l	330431	3	11.644	805237	50.0
cis-1-Ethyl-3-m...	12.144	31.3	ug/l	504004	3	11.644	805237	50.0
m-Menthane, (1S...	12.950	20.3	ug/l	374193	4	13.567	923990	50.0
Pentalene, octa...	13.185	20.3	ug/l	375854	4	13.567	923990	50.0
Cyclohexane, (1...	13.455	11.6	ug/l	214801	4	13.567	923990	50.0
Naphthalene, de...	13.891	21.8	ug/l	401903	4	13.567	923990	50.0
Naphthalene, de...	14.402	13.9	ug/l	257746	4	13.567	923990	50.0
Naphthalene, 1-...	16.685	12.1	ug/l	222855	4	13.567	923990	50.0