

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD120321\
 Data File : VD071133.D
 Acq On : 03 Dec 2021 14:03
 Operator : VA/SY
 Sample : M4920-01
 Misc : 5.00G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 FDR-20211202-WC-S

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: VD071133.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.068	22	25	27	rBV2	1911	2249	1.62%	0.151%
2	2.232	43	53	55	rBV4	6068	14214	10.25%	0.954%
3	2.279	59	61	63	rVV2	2640	2196	1.58%	0.147%
4	2.303	63	65	68	rVV3	2295	2853	2.06%	0.192%
5	2.332	68	70	72	rVV3	2183	2475	1.78%	0.166%
6	2.373	74	77	79	rBV3	2220	2750	1.98%	0.185%
7	2.432	84	87	90	rBV2	2751	3743	2.70%	0.251%
8	2.520	100	102	105	rBV3	2025	1909	1.38%	0.128%
9	2.562	107	109	114	rBV3	2552	4181	3.01%	0.281%
10	2.720	134	136	138	rVB	2603	1753	1.26%	0.118%
11	2.750	138	141	142	rBV3	2768	3069	2.21%	0.206%
12	2.785	144	147	148	rBV2	2959	3187	2.30%	0.214%
13	2.926	168	171	176	rVB3	2078	3209	2.31%	0.215%
14	3.067	192	195	198	rVB2	1214	1669	1.20%	0.112%
15	3.156	207	210	212	rVB	1803	1847	1.33%	0.124%
16	3.209	216	219	222	rVB2	1419	1663	1.20%	0.112%
17	3.532	270	274	275	rBV2	1187	1652	1.19%	0.111%
18	4.156	372	380	382	rBV3	1458	2493	1.80%	0.167%
19	4.285	398	402	405	rBV2	1526	2193	1.58%	0.147%
20	4.426	421	426	430	rBV2	1776	3319	2.39%	0.223%
21	4.791	485	488	491	rBV	1330	2009	1.45%	0.135%
22	4.961	514	517	520	rBV2	1200	1666	1.20%	0.112%
23	5.209	557	559	563	rVV	1322	1895	1.37%	0.127%
24	6.403	759	762	766	rBV2	1394	1769	1.28%	0.119%
25	7.085	875	878	881	rBV2	1311	1959	1.41%	0.131%
26	7.185	888	895	900	rBV3	3482	9788	7.06%	0.657%
27	7.350	921	923	928	rBV2	1009	1875	1.35%	0.126%
28	7.914	1011	1019	1024	rVV3	13935	34337	24.76%	2.305%
29	7.979	1024	1030	1039	rVV2	27316	60201	43.40%	4.041%
30	8.332	1081	1090	1098	rVV3	20354	48679	35.10%	3.268%
31	8.726	1155	1157	1163	rVB3	1645	2616	1.89%	0.176%
32	8.861	1169	1180	1188	rVV2	39239	80426	57.99%	5.399%
33	9.926	1357	1361	1365	rVB2	1114	1654	1.19%	0.111%
34	10.085	1387	1388	1393	rVB2	1370	1692	1.22%	0.114%
35	10.132	1393	1396	1400	rVB2	1968	2128	1.53%	0.143%
36	10.185	1400	1405	1407	rBV2	1088	1888	1.36%	0.127%

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37	10.338	1422	1431	1439	rBV	62440	113585	81.89%	7.624%
38	10.526	1460	1463	1466	rVV2	1660	2449	1.77%	0.164%
39	10.596	1472	1475	1482	rVB	1369	1835	1.32%	0.123%
40	11.202	1575	1578	1581	rBV	1389	1867	1.35%	0.125%
41	11.638	1645	1652	1660	rBV	65842	115736	83.44%	7.769%
42	12.008	1711	1715	1718	rBV2	1634	1796	1.29%	0.121%
43	12.091	1724	1729	1731	rBV2	1702	2618	1.89%	0.176%
44	12.496	1795	1798	1799	rBV2	1438	1740	1.25%	0.117%
45	12.626	1814	1820	1826	rBV2	49777	84388	60.84%	5.664%
46	12.920	1866	1870	1871	rBV2	1800	2182	1.57%	0.146%
47	13.102	1899	1901	1904	rVV3	2019	1687	1.22%	0.113%
48	13.214	1913	1920	1926	rVB2	43778	70959	51.16%	4.763%
49	13.314	1931	1937	1939	rBV2	4230	5356	3.86%	0.360%
50	13.361	1944	1945	1947	rBV2	2157	1948	1.40%	0.131%
51	13.461	1959	1962	1964	rVV3	2345	2380	1.72%	0.160%
52	13.490	1964	1967	1972	rVB5	4731	5790	4.17%	0.389%
53	13.567	1972	1980	1987	rBV	66799	115236	83.08%	7.735%
54	13.679	1998	1999	2003	rBV4	2231	2577	1.86%	0.173%
55	13.726	2003	2007	2008	rBV3	3774	4186	3.02%	0.281%
56	13.790	2016	2018	2019	rBV	2046	1920	1.38%	0.129%
57	13.849	2025	2028	2032	rVB5	3944	6013	4.34%	0.404%
58	13.920	2034	2040	2043	rBV5	11281	23619	17.03%	1.585%
59	13.985	2050	2051	2053	rBV2	3522	2845	2.05%	0.191%
60	14.120	2073	2074	2076	rBV2	3757	2215	1.60%	0.149%
61	14.208	2084	2089	2096	rBV2	81671	138700	100.00%	9.310%
62	14.255	2096	2097	2100	rVB3	5225	5165	3.72%	0.347%
63	14.314	2103	2107	2108	rBV4	2631	3111	2.24%	0.209%
64	14.385	2115	2119	2120	rBV3	3344	4244	3.06%	0.285%
65	14.408	2120	2123	2124	rVV2	2062	2400	1.73%	0.161%
66	14.485	2133	2136	2137	rBV2	2185	2035	1.47%	0.137%
67	14.508	2137	2140	2142	rVB4	4246	3691	2.66%	0.248%
68	14.532	2142	2144	2150	rVB7	3154	5567	4.01%	0.374%
69	14.667	2165	2167	2168	rVV2	3033	2263	1.63%	0.152%
70	14.685	2168	2170	2173	rVB4	3615	3994	2.88%	0.268%
71	14.832	2192	2195	2196	rBV3	2085	2042	1.47%	0.137%
72	14.849	2196	2198	2199	rVV2	5982	2846	2.05%	0.191%
73	14.890	2199	2205	2210	rVV	50363	87109	62.80%	5.847%
74	14.949	2214	2215	2217	rVB2	2849	1734	1.25%	0.116%
75	14.979	2217	2220	2223	rBV5	5257	6960	5.02%	0.467%
76	15.067	2231	2235	2236	rBV4	11801	13722	9.89%	0.921%

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77	15.196	2253	2257	2259	rBV4	3217	5463	3.94%	0.367%
78	15.302	2273	2275	2278	rBV4	2116	2676	1.93%	0.180%
79	15.355	2278	2284	2290	rBV2	25901	44802	32.30%	3.007%
80	15.426	2290	2296	2301	rBV8	6021	12063	8.70%	0.810%
81	15.549	2315	2317	2318	rBV2	3336	1904	1.37%	0.128%
82	15.602	2324	2326	2328	rBV3	3481	3198	2.31%	0.215%
83	15.643	2329	2333	2334	rBV3	16039	17765	12.81%	1.192%
84	15.684	2336	2340	2343	rVB2	14760	24968	18.00%	1.676%
85	15.755	2350	2352	2355	rVB4	11550	11418	8.23%	0.766%
86	15.820	2361	2363	2364	rBV2	3990	2718	1.96%	0.182%
87	15.873	2371	2372	2373	rBV	4841	2026	1.46%	0.136%
88	15.943	2383	2384	2385	rBV	5391	3607	2.60%	0.242%
89	16.014	2394	2396	2397	rVB2	3876	1750	1.26%	0.117%
90	16.037	2397	2400	2406	rVB8	7229	14868	10.72%	0.998%
91	16.173	2418	2423	2431	rVB4	24499	49151	35.44%	3.299%
92	16.296	2440	2444	2448	rVB7	9698	18434	13.29%	1.237%
93	16.331	2448	2450	2451	rBV2	3299	2726	1.97%	0.183%
94	16.349	2451	2453	2454	rBV2	4000	2665	1.92%	0.179%
95	16.414	2462	2464	2465	rBV2	4699	4591	3.31%	0.308%
96	16.449	2465	2470	2479	rVB10	13579	40721	29.36%	2.733%
97	16.520	2479	2482	2484	rBV4	3665	4522	3.26%	0.304%
98	16.573	2486	2491	2494	rBV7	6530	12399	8.94%	0.832%
99	16.620	2494	2499	2504	rVB4	11693	23625	17.03%	1.586%
100	16.761	2521	2523	2525	rVB3	4595	3995	2.88%	0.268%

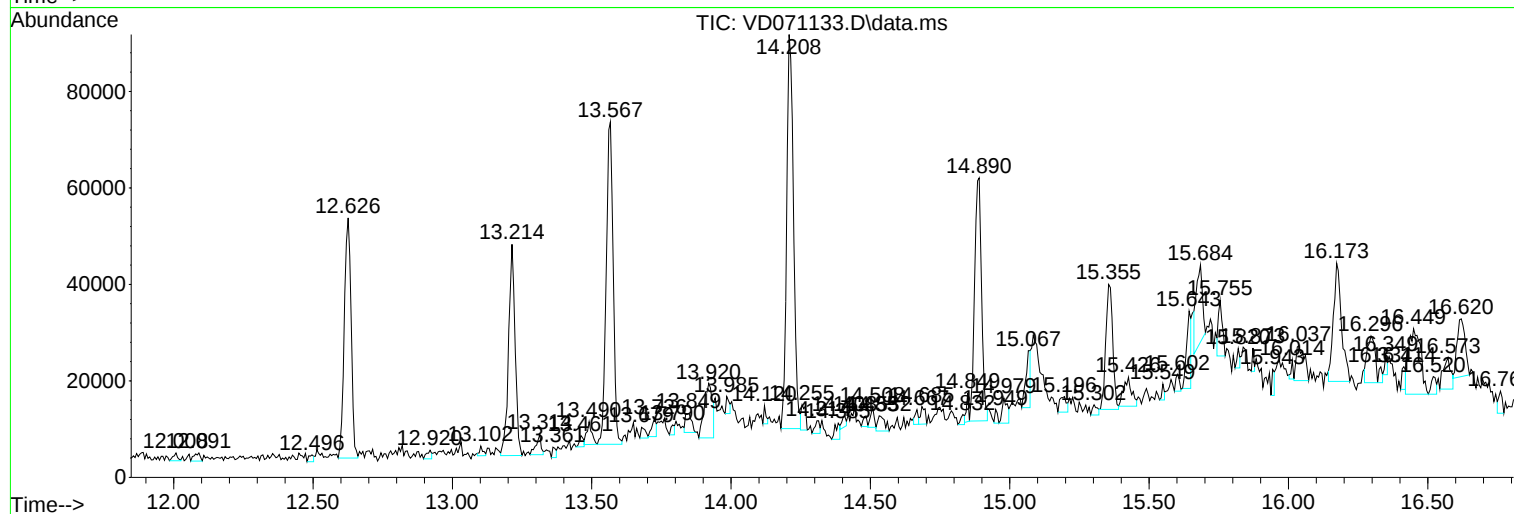
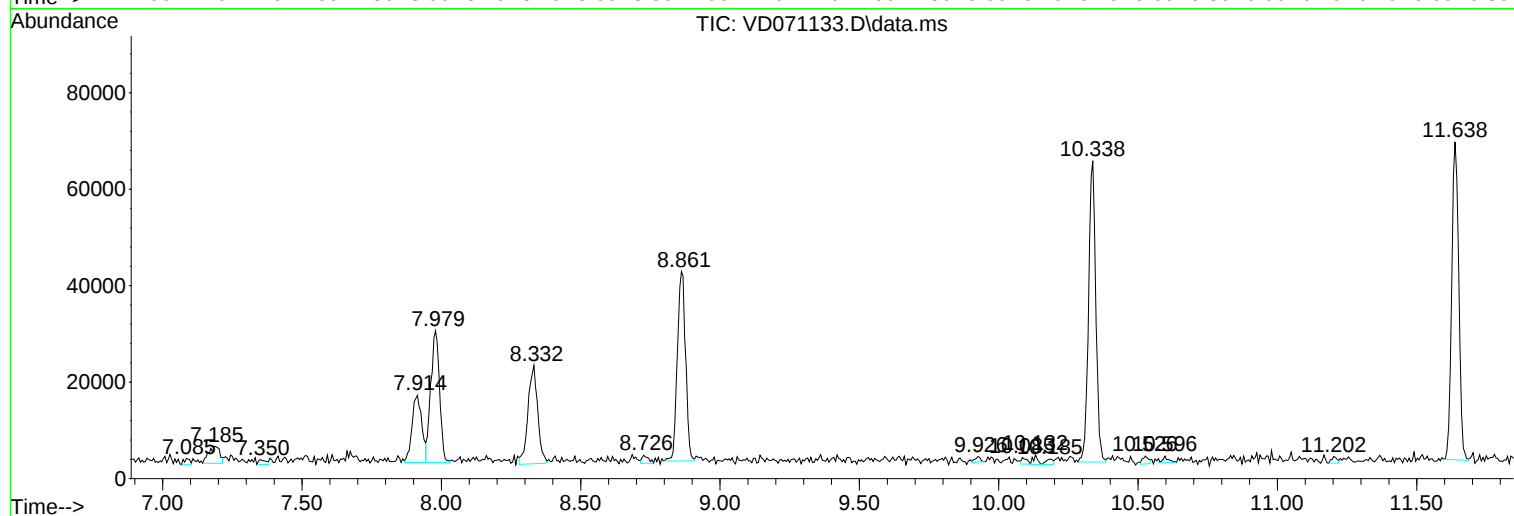
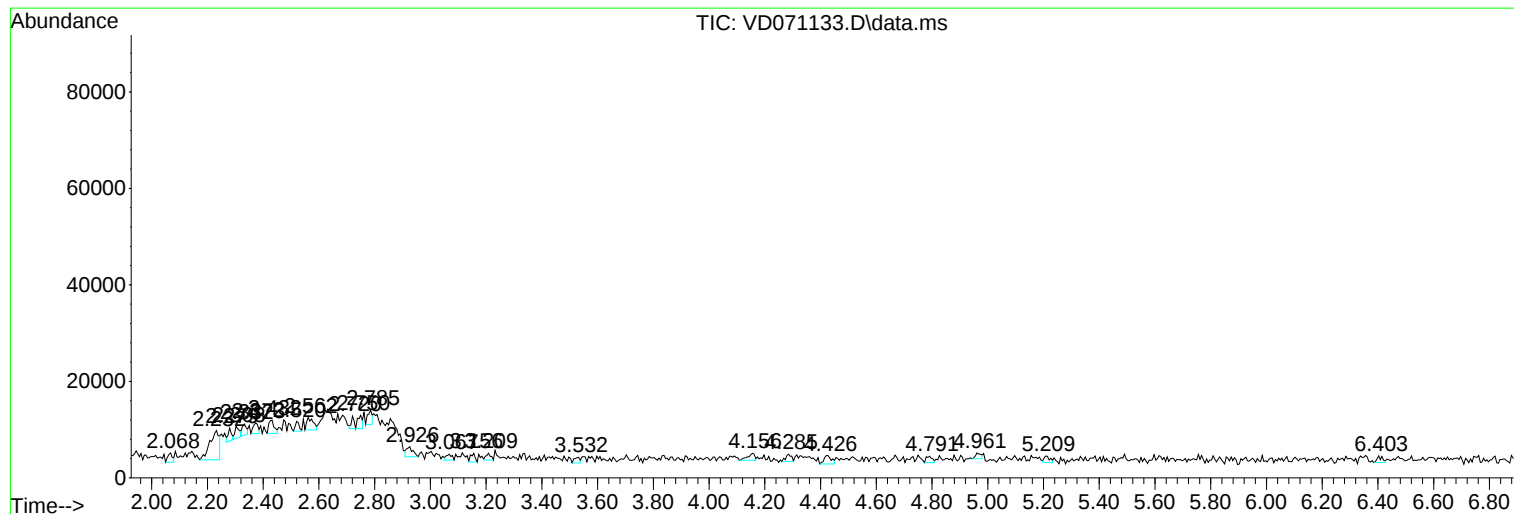
Sum of corrected areas: 1489771

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Instrument :
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ClientSampleId :
FDR-20211202-WC-S

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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FDR-20211202-WC-S

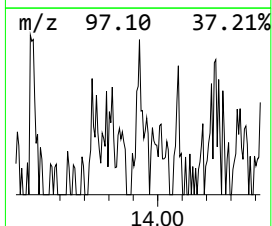
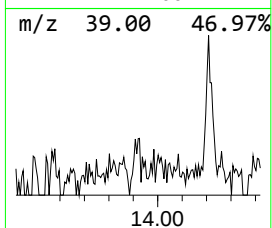
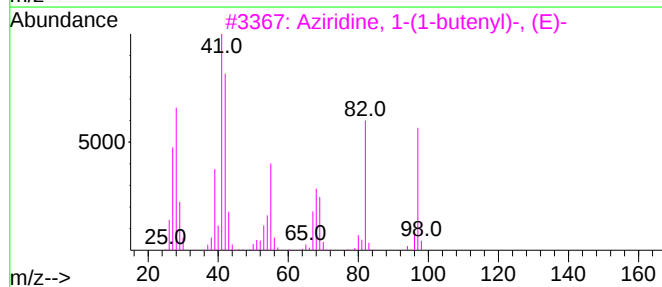
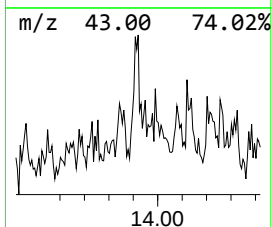
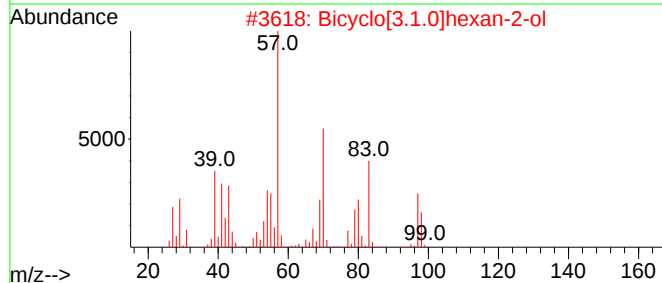
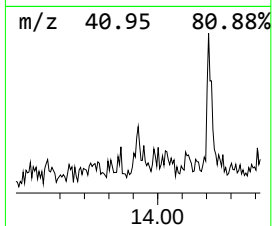
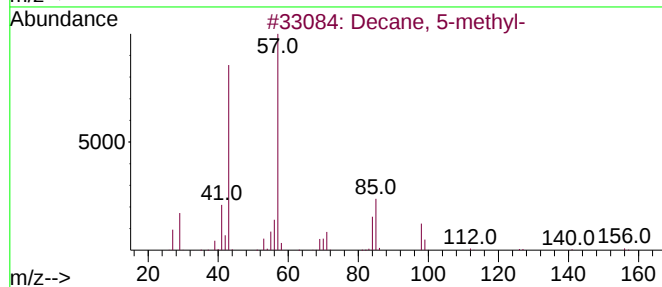
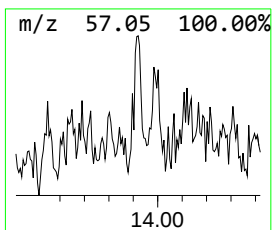
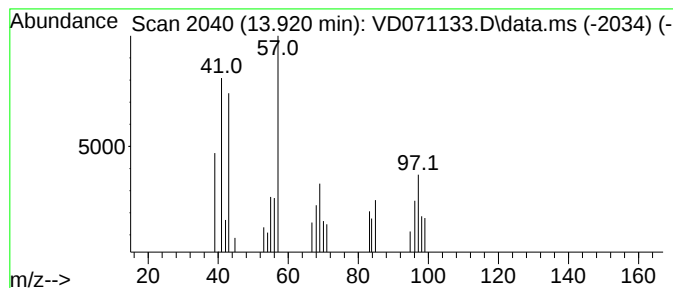
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

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

Peak Number 2 unknown13.920 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.920	10.25 ug/l	23619	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 5-methyl-	156	C11H24	013151-35-4	27
2			Bicyclo[3.1.0]hexan-2-ol	98	C6H10O	1000194-18-6	22
3			Aziridine, 1-(1-butenyl)-, (E)-	97	C6H11N	080839-92-5	18
4			Nonane, 2,5-dimethyl-	156	C11H24	017302-27-1	16
5			Decane, 2,5,6-trimethyl-	184	C13H28	062108-23-0	16



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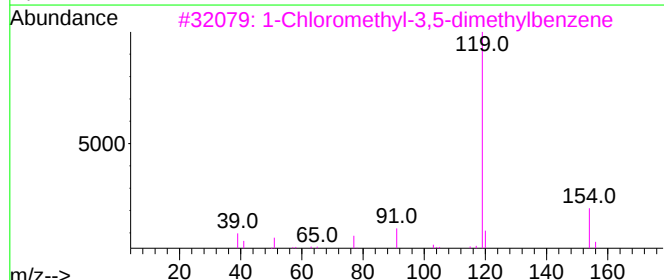
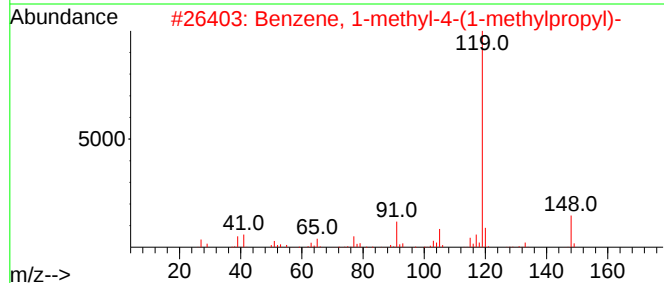
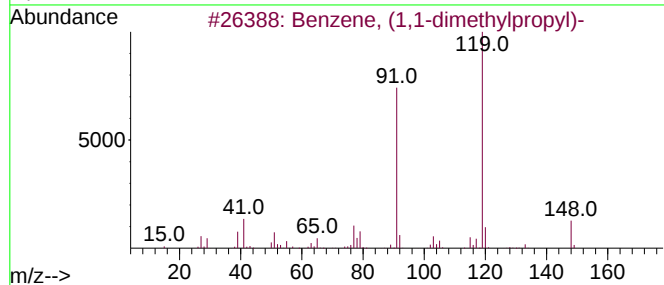
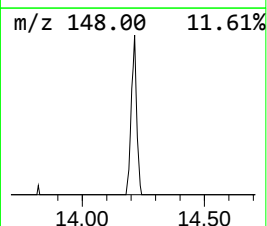
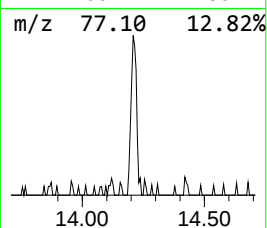
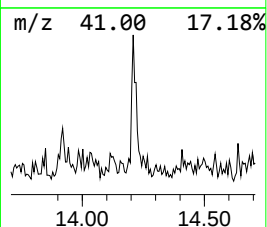
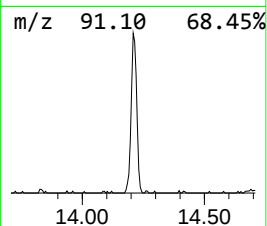
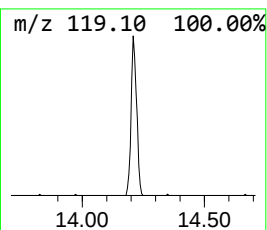
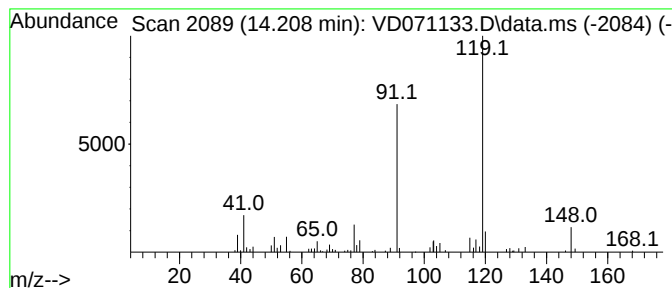
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, (1,1-dimethylpropyl)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.208	60.18 ug/l	138700	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	94
2			Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	87
3			1-Chloromethyl-3,5-dimethylbenzene	154	C9H11Cl	002745-54-2	72
4			Benzene, tert-butyl-	134	C10H14	000098-06-6	72
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	72



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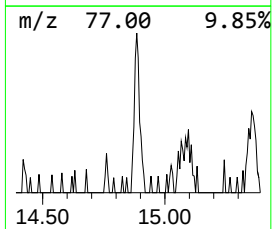
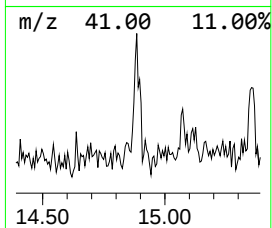
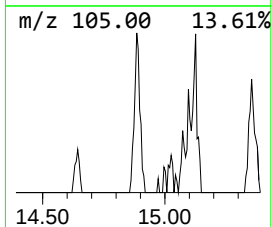
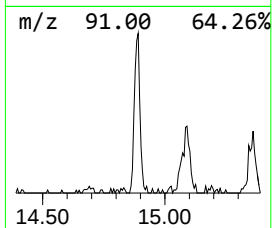
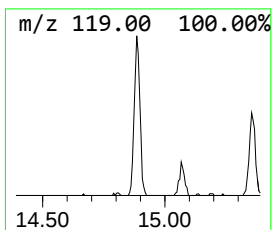
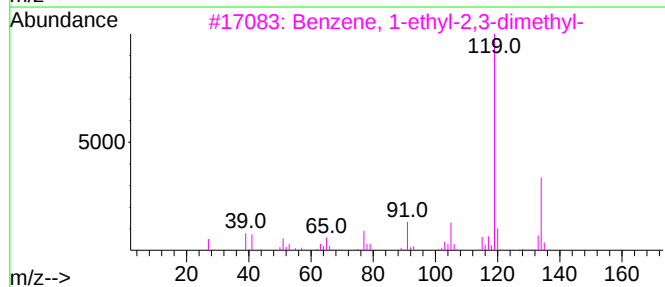
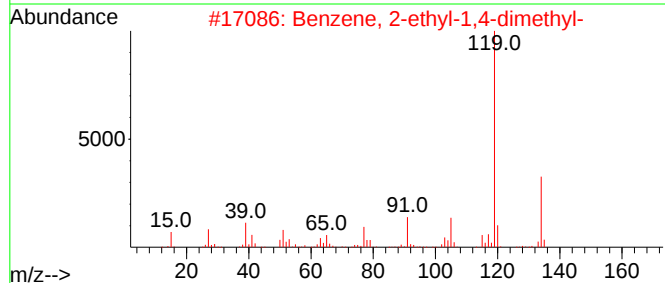
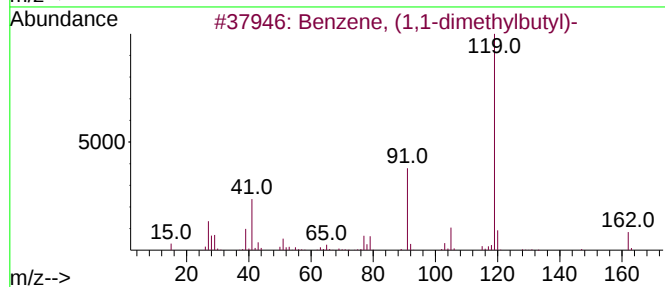
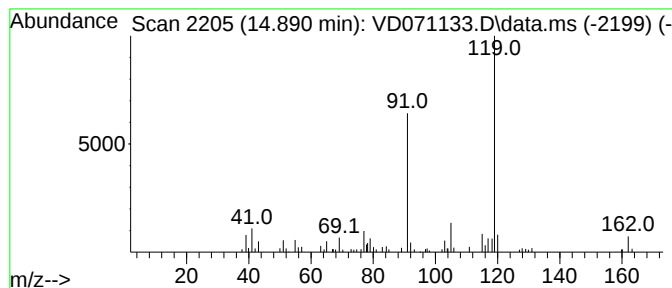
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, (1,1-dimethylbutyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.890	37.80 ug/l	87109	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1,1-dimethylbutyl)-	162	C12H18	001985-57-5	80
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	78
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	78
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	78
5			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD120321\
Data File : VD071133.D
Acq On : 03 Dec 2021 14:03
Operator : VA/SY
Sample : M4920-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
FDR-20211202-WC-S

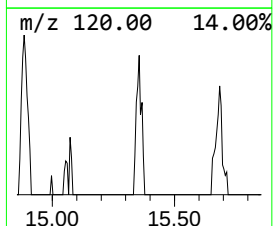
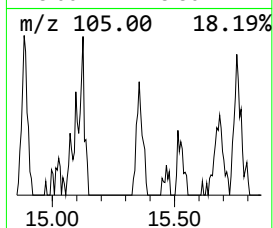
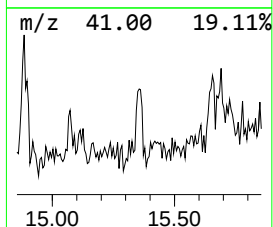
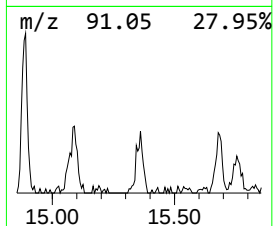
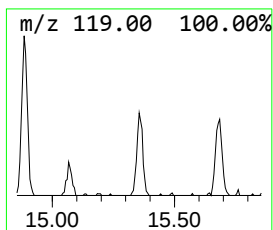
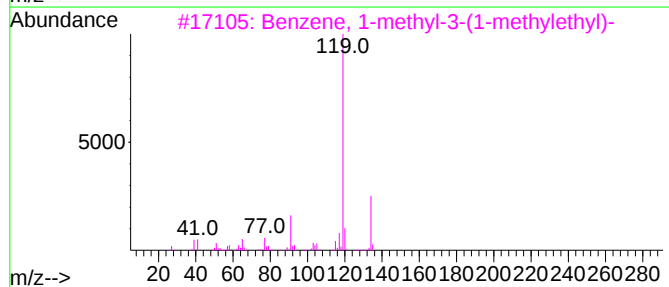
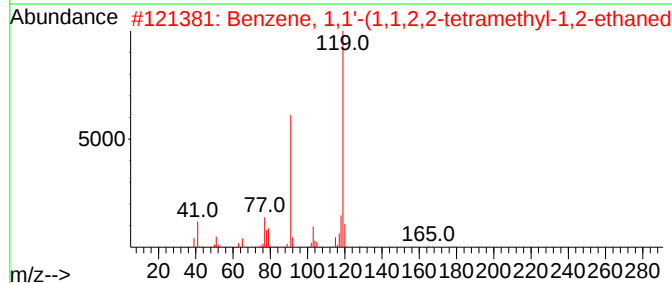
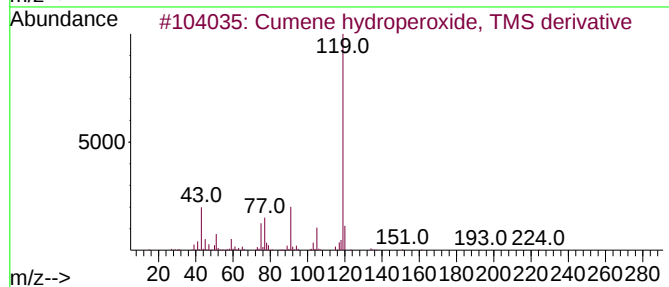
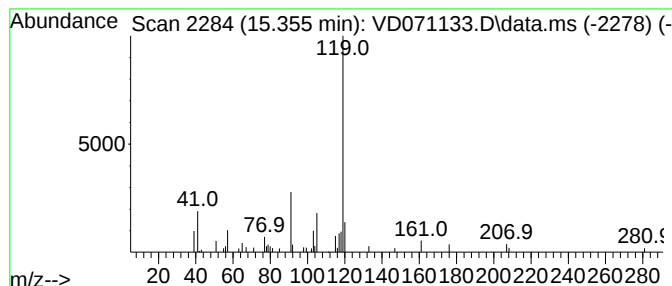
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Quant Title : SW846 8260

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

Peak Number 5 Cumene hydroperoxide, TMS d... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.355	19.44 ug/l	44802	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cumene hydroperoxide, TMS deriva...	224	C12H20O2Si	018057-16-4	72
2			Benzene, 1,1'-(1,1,2,2-tetrameth...	238	C18H22	001889-67-4	64
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	59
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	59
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD120321\
Data File : VD071133.D
Acq On : 03 Dec 2021 14:03
Operator : VA/SY
Sample : M4920-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
FDR-20211202-WC-S

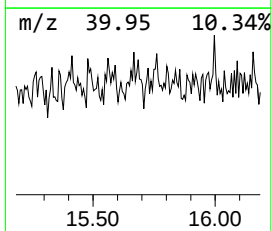
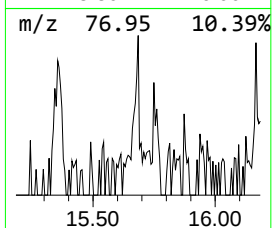
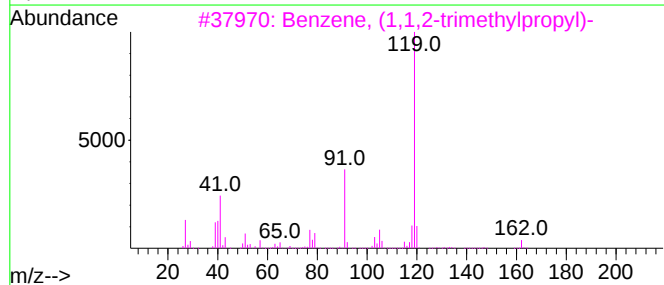
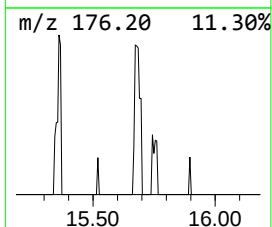
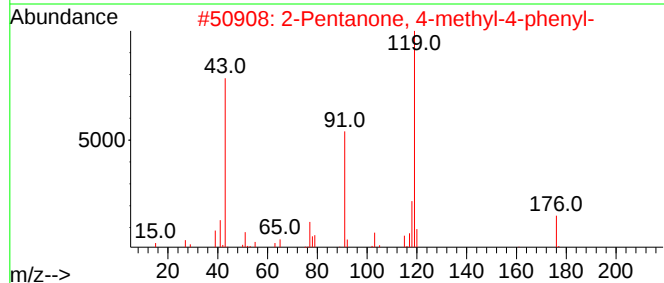
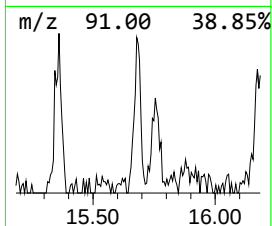
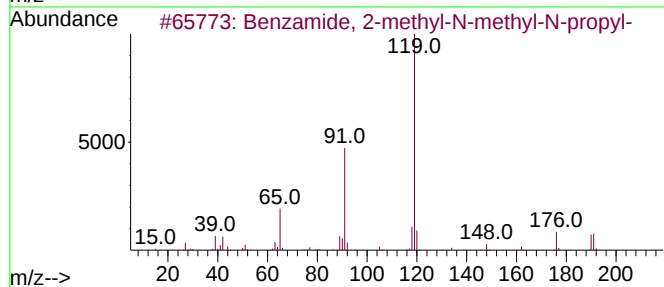
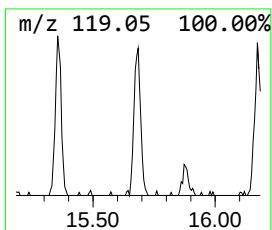
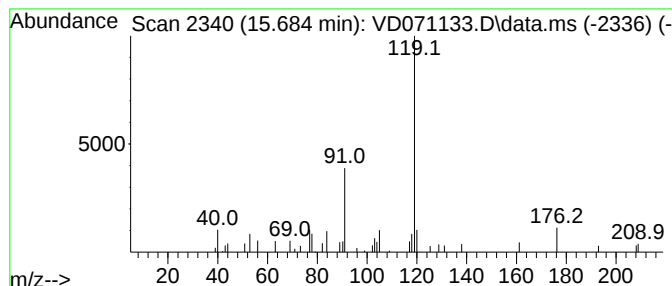
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Quant Title : SW846 8260

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

Peak Number 6 Benzamide, 2-methyl-N-methy... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.685	10.83 ug/l	24968	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzamide, 2-methyl-N-methyl-N-p...	191	C12H17NO	1000421-44-4	59
2			2-Pentanone, 4-methyl-4-phenyl-	176	C12H16O	007403-42-1	58
3			Benzene, (1,1,2-trimethylpropyl)-	162	C12H18	026356-11-6	53
4			Benzene, 1,1'-(1,1,2,2-tetrameth...	238	C18H22	001889-67-4	53
5			N-(m-Toluoyl)glycine	193	C10H11NO3	027115-49-7	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD120321\
Data File : VD071133.D
Acq On : 03 Dec 2021 14:03
Operator : VA/SY
Sample : M4920-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
FDR-20211202-WC-S

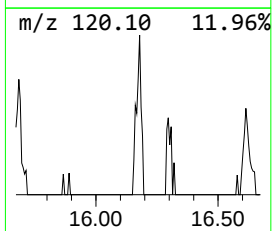
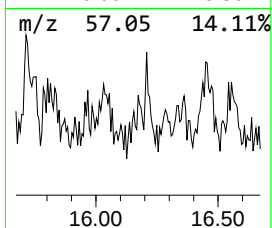
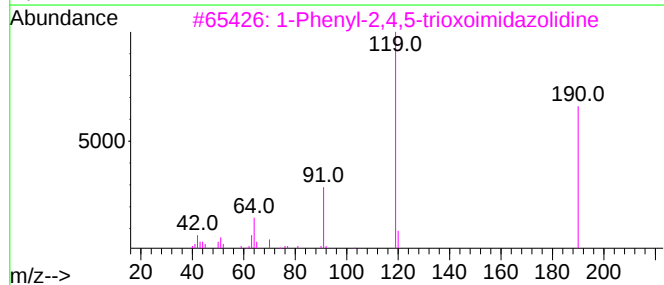
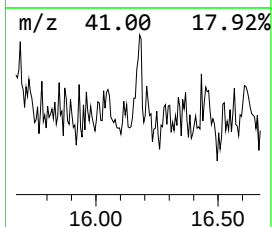
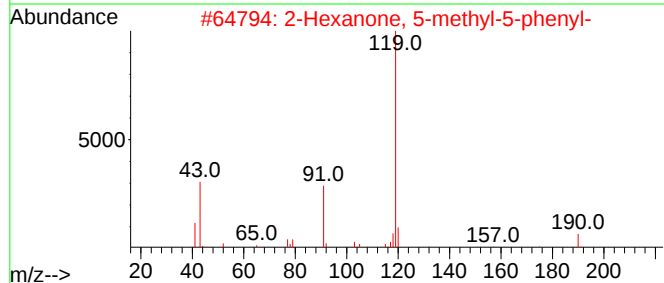
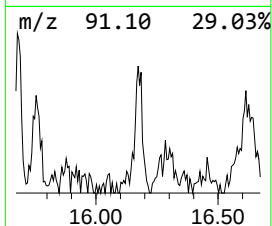
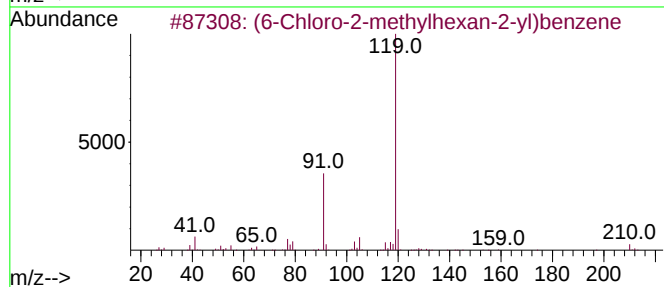
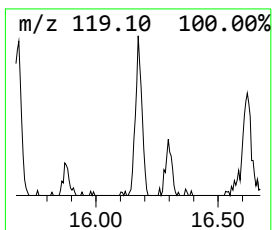
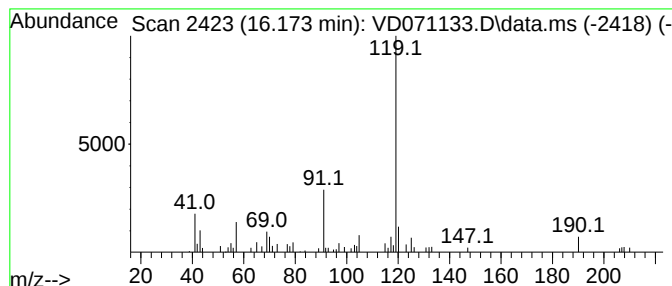
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Quant Title : SW846 8260

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

Peak Number 7 (6-Chloro-2-methylhexan-2-yl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.173	21.33 ug/l	49151	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(6-Chloro-2-methylhexan-2-yl)ben...	210	C13H19Cl	1417621-45-4	81
2			2-Hexanone, 5-methyl-5-phenyl-	190	C13H18O	014128-61-1	72
3			1-Phenyl-2,4,5-trioxoimidazolidine	190	C9H6N2O3	002211-33-8	64
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	64
5			Benzene, 1-ethyl-4-(2-methylprop...	162	C12H18	100319-40-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD120321\
Data File : VD071133.D
Acq On : 03 Dec 2021 14:03
Operator : VA/SY
Sample : M4920-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
FDR-20211202-WC-S

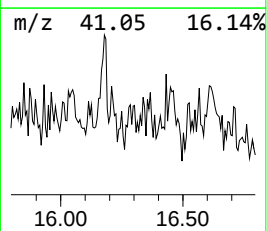
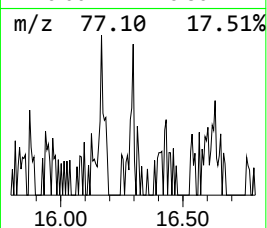
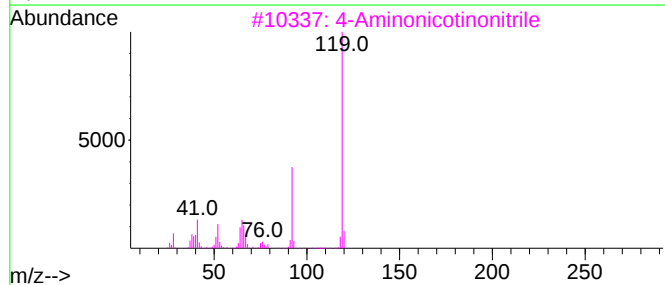
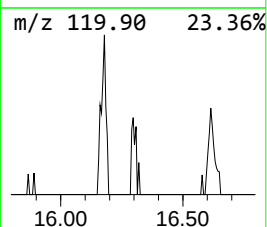
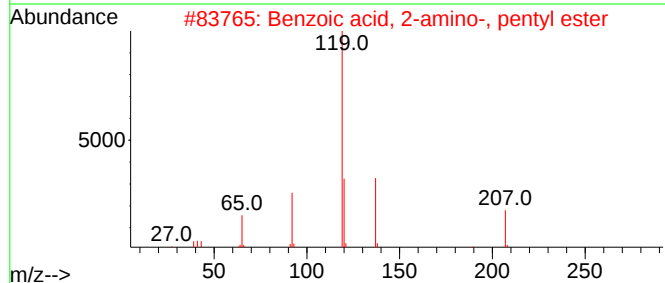
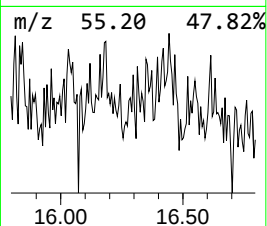
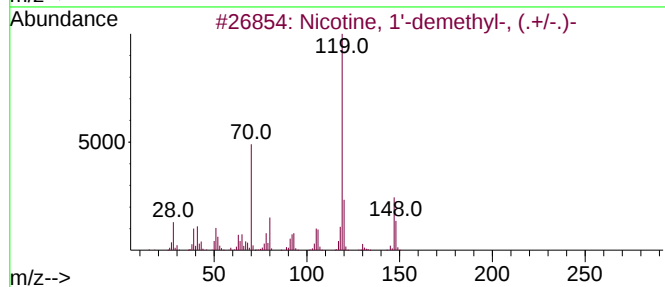
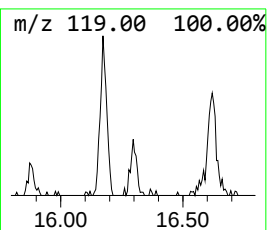
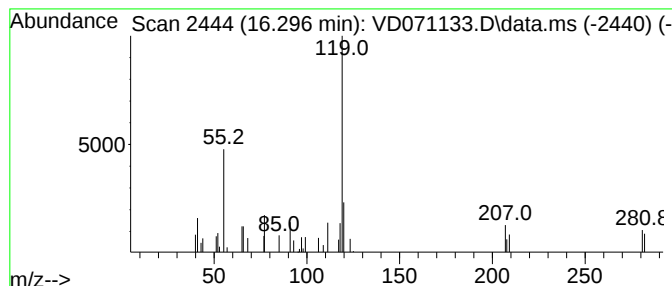
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Quant Title : SW846 8260

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

Peak Number 8 unknown16.296 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.296	8.00 ug/l	18434	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nicotine, 1'-demethyl-, (+/-)-	148	C9H12N2	005746-86-1	47
2			Benzoic acid, 2-amino-, pentyl e...	207	C12H17NO2	030100-15-3	47
3			4-Aminonicotinonitrile	119	C6H5N3	015827-84-6	43
4			5-Aza-2H-indazole	119	C6H5N3	000271-50-1	37
5			Carbamate, N-phenyl-, (3-methyl-...	207	C11H13NO3	137531-94-3	33



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD120321\
Data File : VD071133.D
Acq On : 03 Dec 2021 14:03
Operator : VA/SY
Sample : M4920-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
FDR-20211202-WC-S

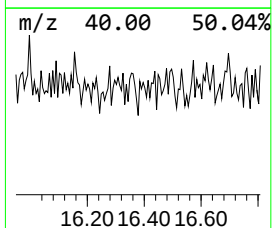
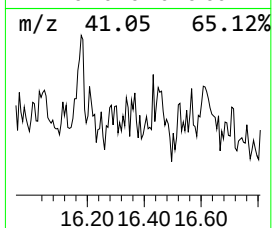
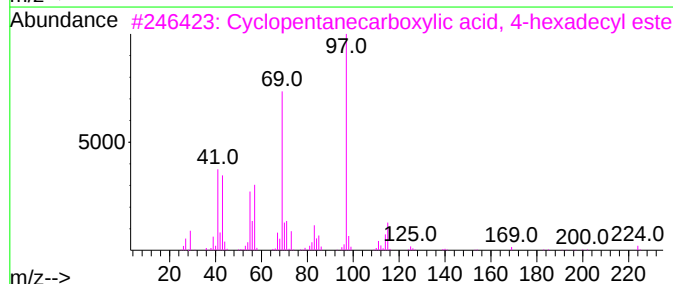
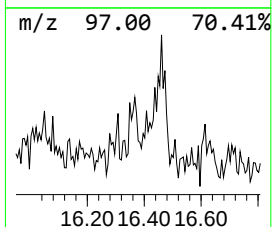
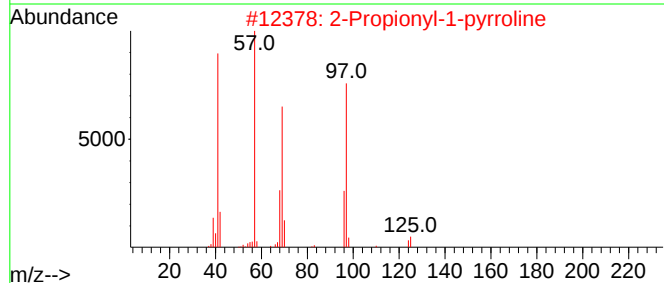
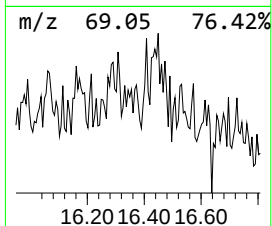
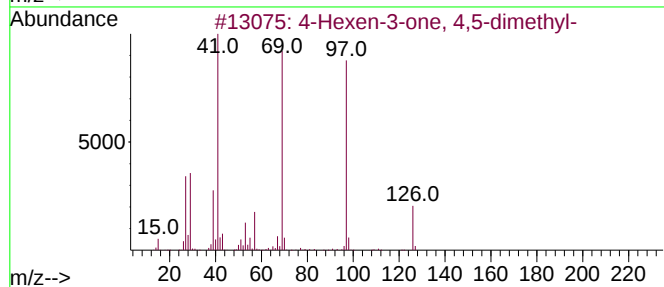
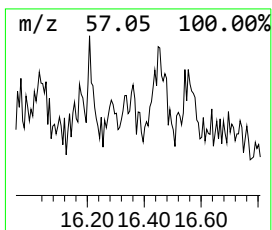
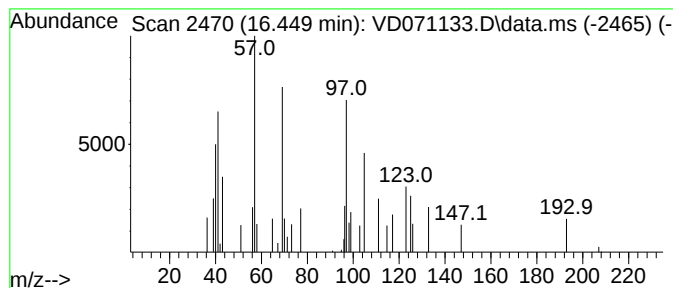
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Quant Title : SW846 8260

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

Peak Number 9 unknown16.449 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.449	17.67 ug/l	40721	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Hexen-3-one, 4,5-dimethyl-	126	C8H14O	017325-90-5	43
2			2-Propionyl-1-pyrroline	125	C7H11NO	1010298-55-4	43
3			Cyclopentanecarboxylic acid, 4-h...	338	C22H42O2	1000282-77-5	32
4			1-Ethyl-2-hydroxymethylimidazole	126	C6H10N2O	063634-44-6	25
5			2-Methyl-2-octene	126	C9H18	016993-86-5	22



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD120321\
Data File : VD071133.D
Acq On : 03 Dec 2021 14:03
Operator : VA/SY
Sample : M4920-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
FDR-20211202-WC-S

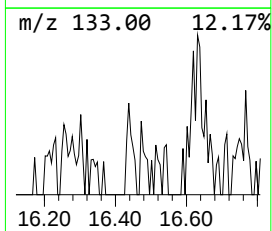
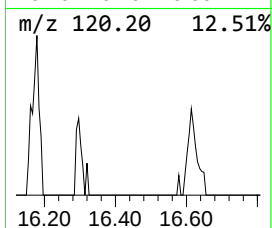
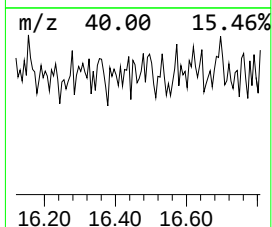
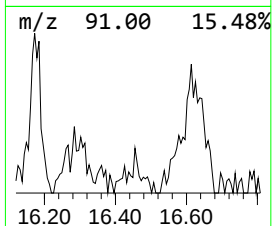
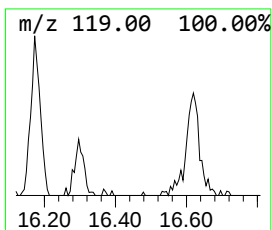
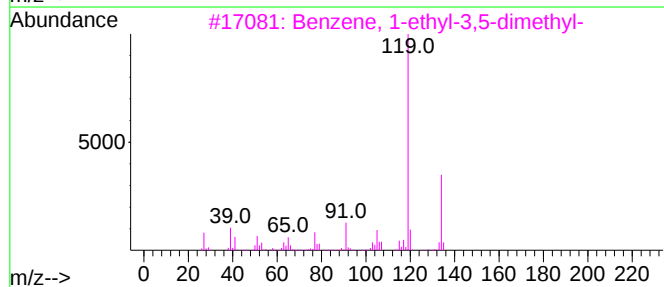
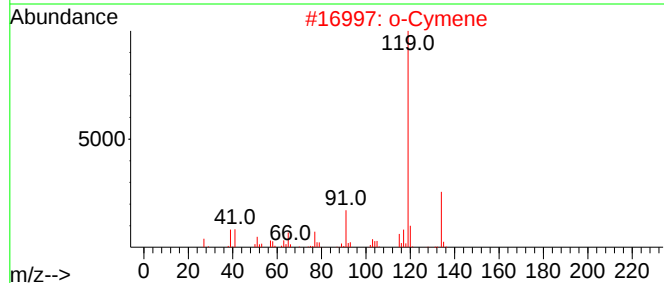
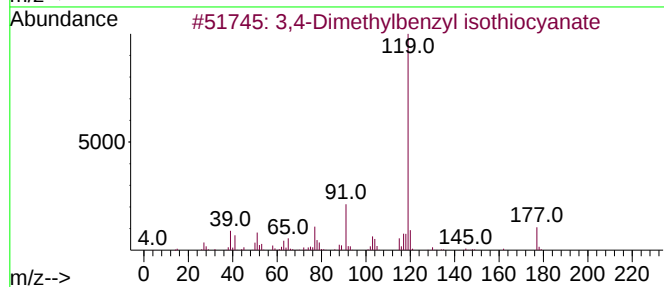
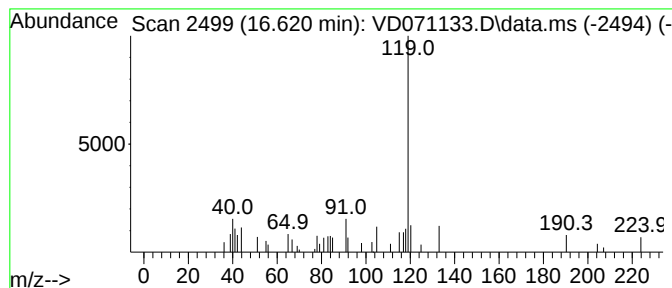
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Quant Title : SW846 8260

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

Peak Number 10 3,4-Dimethylbenzyl isothioc... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.620	10.25 ug/l	23625	1,4-Dichlorobenzene-d4	13.567

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3,4-Dimethylbenzyl isothiocyanate	177	C10H11NS	206559-59-3	64
2		o-Cymene	134	C10H14	000527-84-4	59
3		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	59
4		Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	59
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD120321\
Data File : VD071133.D
Acq On : 03 Dec 2021 14:03
Operator : VA/SY
Sample : M4920-01
Misc : 5.00G/5.00ml/MSVOA_D/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
FDR-20211202-WC-S

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.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
unknown13.920	13.920	10.3	ug/l	23619	4	13.567	115236	50.0
Benzene, (1,1-d...	14.208	60.2	ug/l	138700	4	13.567	115236	50.0
Benzene, (1,1-d...	14.890	37.8	ug/l	87109	4	13.567	115236	50.0
Cumene hydroper...	15.355	19.4	ug/l	44802	4	13.567	115236	50.0
Benzamide, 2-me...	15.685	10.8	ug/l	24968	4	13.567	115236	50.0
(6-Chloro-2-met...	16.173	21.3	ug/l	49151	4	13.567	115236	50.0
unknown16.296	16.296	8.0	ug/l	18434	4	13.567	115236	50.0
unknown16.449	16.449	17.7	ug/l	40721	4	13.567	115236	50.0
3,4-Dimethylben...	16.620	10.3	ug/l	23625	4	13.567	115236	50.0