

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070622\
 Data File : VN073345.D
 Acq On : 06 Jul 2022 15:11
 Operator : JC\MD
 Sample : N3564-04
 Misc : 4.97g/5mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-14B-11-1-GR

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_N\methods\82N062822W.M
 Title : SW846 8260

Signal : TIC: VN073345.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.946	176	180	187	rBV2	65914	103547	4.87%	0.304%
2	3.322	236	244	254	rVB4	33221	82648	3.89%	0.242%
3	3.834	326	331	339	rVB3	12318	29899	1.41%	0.088%
4	4.775	482	491	502	rBV2	17427	54165	2.55%	0.159%
5	4.922	504	516	519	rVV3	50347	153495	7.22%	0.450%
6	4.987	520	527	543	rVB3	117694	416601	19.59%	1.221%
7	5.210	555	565	571	rBV2	16569	55609	2.62%	0.163%
8	5.475	597	610	629	rBV2	249786	877572	41.27%	2.572%
9	6.346	748	758	769	rVB2	51287	148455	6.98%	0.435%
10	6.775	819	831	843	rBV6	35275	121121	5.70%	0.355%
11	6.928	848	857	867	rBV3	41056	127378	5.99%	0.373%
12	7.034	867	875	882	rBV2	14259	35583	1.67%	0.104%
13	7.287	911	918	929	rVB2	16822	47143	2.22%	0.138%
14	7.751	990	997	1007	rVB4	22839	65483	3.08%	0.192%
15	7.945	1017	1030	1033	rBV	193574	473456	22.27%	1.388%
16	8.004	1033	1040	1048	rVV3	571381	1553733	73.08%	4.554%
17	8.087	1048	1054	1065	rVB	190892	471370	22.17%	1.382%
18	8.210	1065	1075	1081	rBV2	219186	526539	24.76%	1.543%
19	8.269	1081	1085	1093	rVB2	134295	291016	13.69%	0.853%
20	8.363	1093	1101	1111	rVB	198011	450148	21.17%	1.320%
21	8.487	1111	1122	1128	rBV2	258625	633336	29.79%	1.857%
22	8.557	1128	1134	1139	rVV2	200022	430977	20.27%	1.263%
23	8.628	1139	1146	1157	rVB	451414	1043636	49.08%	3.059%
24	8.898	1180	1192	1205	rBV	596302	1436097	67.54%	4.210%
25	9.051	1211	1218	1223	rVB2	18198	34492	1.62%	0.101%
26	9.128	1223	1231	1235	rBV2	22668	45755	2.15%	0.134%
27	9.198	1235	1243	1256	rVB2	22183	62237	2.93%	0.182%
28	9.387	1260	1275	1281	rBV2	733094	2126213	100.00%	6.233%
29	9.439	1281	1284	1294	rVB2	103452	180841	8.51%	0.530%
30	9.569	1299	1306	1311	rVV3	61915	121957	5.74%	0.357%
31	9.657	1311	1321	1330	rVB	244881	538552	25.33%	1.579%
32	9.804	1338	1346	1357	rVB2	221455	447011	21.02%	1.310%
33	9.904	1357	1363	1366	rBV	44754	81345	3.83%	0.238%
34	9.951	1366	1371	1375	rBV2	64108	101127	4.76%	0.296%
35	10.004	1375	1380	1385	rBV4	89172	171447	8.06%	0.503%
36	10.139	1394	1403	1408	rBV4	146720	379350	17.84%	1.112%

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Integrator: RTE

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
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37	10.257	1416	1423	1427	rBV3	113222	237196	11.16%	0.695%
38	10.375	1435	1443	1456	rBV2	955641	2063824	97.07%	6.050%
39	10.498	1459	1464	1466	rVV2	37764	57736	2.72%	0.169%
40	10.557	1466	1474	1481	rVB5	186608	511573	24.06%	1.500%
41	10.669	1481	1493	1496	rBV3	78793	164292	7.73%	0.482%
42	10.716	1496	1501	1508	rVB	259106	480302	22.59%	1.408%
43	10.810	1508	1517	1523	rBV4	159698	390218	18.35%	1.144%
44	10.898	1528	1532	1537	rVB2	58819	93933	4.42%	0.275%
45	10.986	1537	1547	1553	rBV	146945	287445	13.52%	0.843%
46	11.086	1553	1564	1572	rBV5	109346	393208	18.49%	1.153%
47	11.157	1572	1576	1579	rVV6	64086	113233	5.33%	0.332%
48	11.192	1579	1582	1585	rVV4	69517	117383	5.52%	0.344%
49	11.228	1585	1588	1592	rVV	104349	163805	7.70%	0.480%
50	11.281	1592	1597	1603	rVV	301905	549005	25.82%	1.609%
51	11.333	1603	1606	1610	rVV2	56382	81264	3.82%	0.238%
52	11.386	1610	1615	1620	rVB3	96815	168474	7.92%	0.494%
53	11.486	1620	1632	1640	rVB6	140830	421196	19.81%	1.235%
54	11.569	1640	1646	1649	rBV	66779	119004	5.60%	0.349%
55	11.680	1659	1665	1678	rBV	762345	1453645	68.37%	4.261%
56	11.816	1684	1688	1692	rVB2	63713	85109	4.00%	0.249%
57	11.869	1692	1697	1702	rBV8	54129	110264	5.19%	0.323%
58	11.928	1702	1707	1712	rBV3	88062	192123	9.04%	0.563%
59	12.033	1719	1725	1727	rBV3	39838	75566	3.55%	0.222%
60	12.080	1731	1733	1738	rVB3	17481	34915	1.64%	0.102%
61	12.192	1744	1752	1759	rBV5	189795	474733	22.33%	1.392%
62	12.310	1764	1772	1776	rVB2	115156	229413	10.79%	0.672%
63	12.357	1776	1780	1781	rBV2	47618	62284	2.93%	0.183%
64	12.398	1781	1787	1788	rVV3	148007	269952	12.70%	0.791%
65	12.422	1789	1791	1797	rVB2	165994	247896	11.66%	0.727%
66	12.669	1827	1833	1842	rBV	678956	1169122	54.99%	3.427%
67	12.757	1843	1848	1853	rVV6	64857	139560	6.56%	0.409%
68	12.833	1856	1861	1868	rVB3	216296	417276	19.63%	1.223%
69	12.916	1868	1875	1880	rBV8	97054	243070	11.43%	0.713%
70	12.992	1882	1888	1894	rVV6	145315	383466	18.04%	1.124%
71	13.051	1894	1898	1903	rVB5	76937	134468	6.32%	0.394%
72	13.139	1908	1913	1916	rBV3	75037	123662	5.82%	0.362%
73	13.175	1917	1919	1923	rVV4	35237	49540	2.33%	0.145%
74	13.227	1923	1928	1936	rVB3	154126	277253	13.04%	0.813%
75	13.316	1940	1943	1946	rBV4	27286	38758	1.82%	0.114%
76	13.351	1946	1949	1953	rVV5	41565	48074	2.26%	0.141%

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77	13.439	1961	1964	1967	rVB3	51711	60682	2.85%	0.178%
78	13.498	1968	1974	1979	rVV7	92756	227776	10.71%	0.668%
79	13.539	1979	1981	1985	rVV5	61094	82936	3.90%	0.243%
80	13.610	1987	1993	2002	rVB	871339	1495534	70.34%	4.384%
81	13.757	2014	2018	2023	rBV4	47689	95090	4.47%	0.279%
82	13.933	2043	2048	2053	rBV3	262417	425627	20.02%	1.248%
83	14.063	2064	2070	2076	rBV10	69522	187924	8.84%	0.551%
84	14.151	2082	2085	2089	rVB2	94497	129218	6.08%	0.379%
85	14.263	2098	2104	2109	rBV3	165274	325946	15.33%	0.955%
86	14.327	2110	2115	2121	rVB7	70483	170938	8.04%	0.501%
87	14.404	2121	2128	2130	rBV3	65581	151869	7.14%	0.445%
88	14.445	2130	2135	2138	rVV3	193588	317014	14.91%	0.929%
89	14.480	2138	2141	2145	rVB3	104021	152878	7.19%	0.448%
90	14.557	2151	2154	2158	rBV	78359	123493	5.81%	0.362%
91	14.610	2158	2163	2170	rVV4	164300	366110	17.22%	1.073%
92	14.792	2191	2194	2198	rVB2	83851	106834	5.02%	0.313%
93	14.851	2199	2204	2210	rBV5	195476	493806	23.22%	1.448%
94	15.127	2247	2251	2255	rBV3	139770	192457	9.05%	0.564%
95	15.239	2265	2270	2276	rVB3	213748	446635	21.01%	1.309%
96	15.327	2281	2285	2291	rVB5	138205	286604	13.48%	0.840%
97	15.398	2292	2297	2306	rBV6	185026	377742	17.77%	1.107%
98	15.727	2346	2353	2354	rBV6	100905	195010	9.17%	0.572%
99	16.251	2436	2442	2448	rBV5	205310	487012	22.91%	1.428%
100	16.727	2517	2523	2533	rBV8	211529	622657	29.28%	1.825%

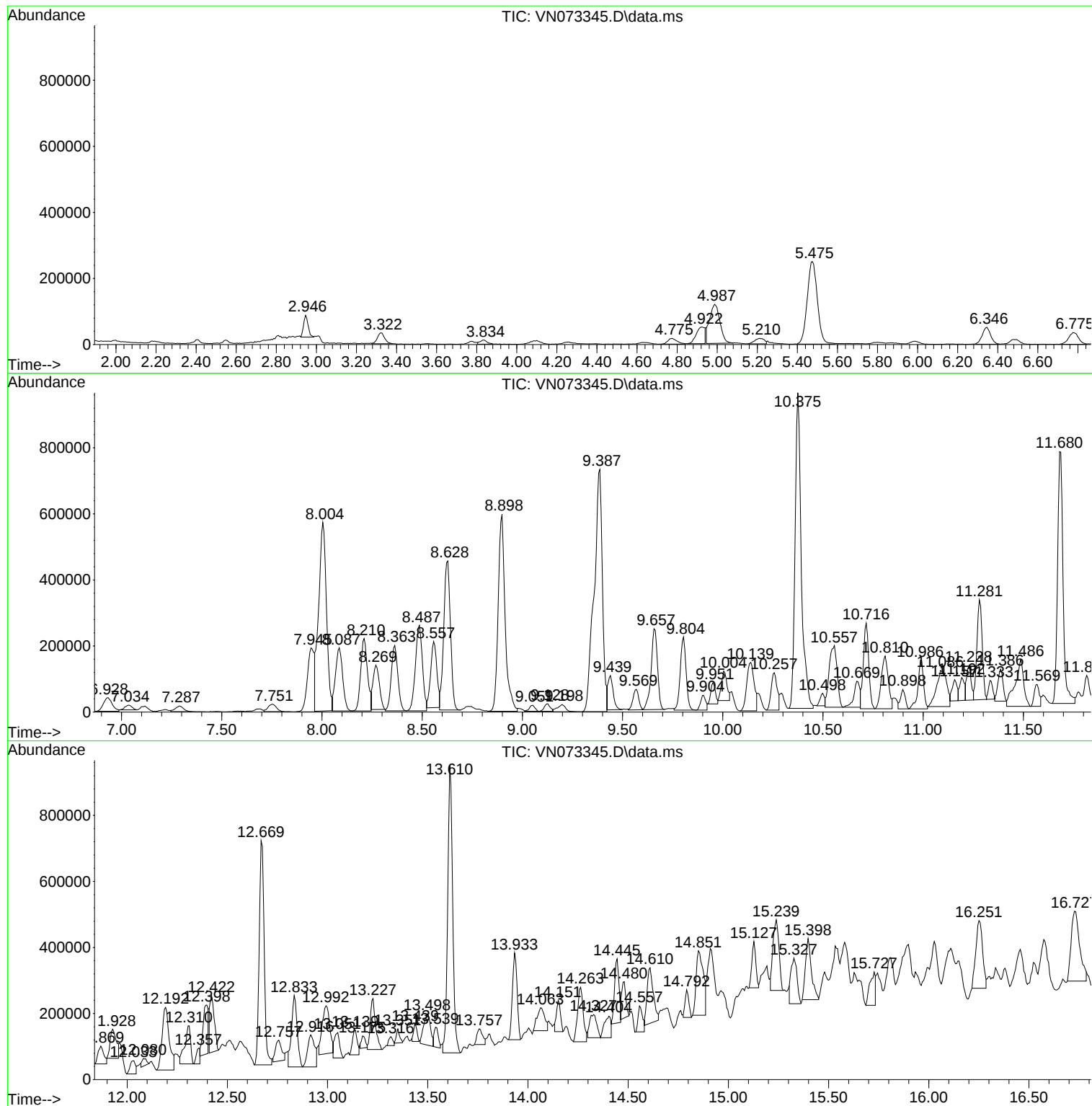
Sum of corrected areas: 34114366

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
Quant Title : SW846 8260

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



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WC-14B-11-1-GR

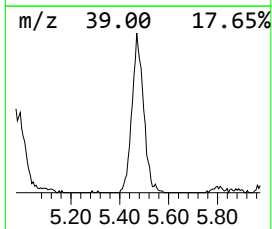
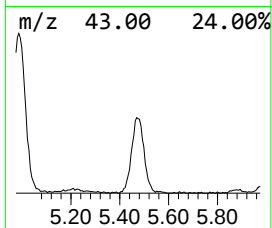
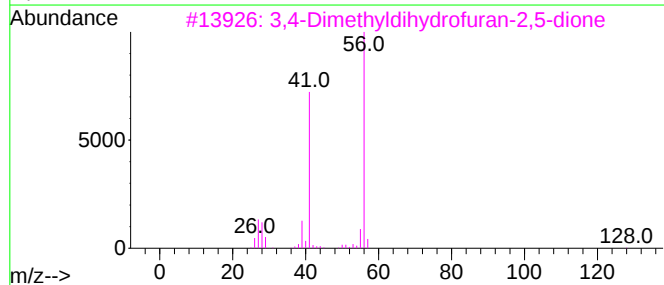
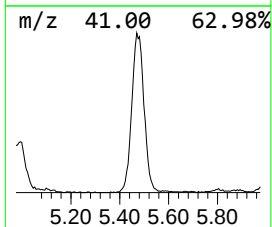
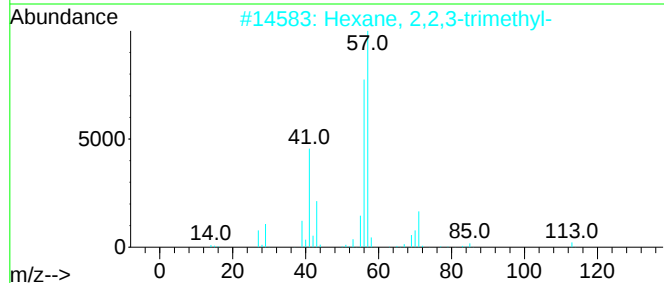
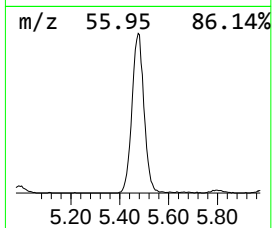
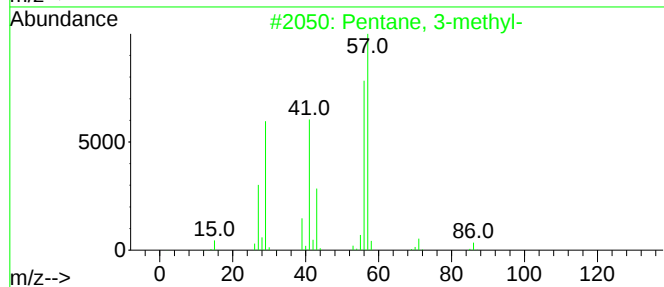
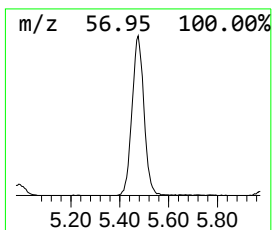
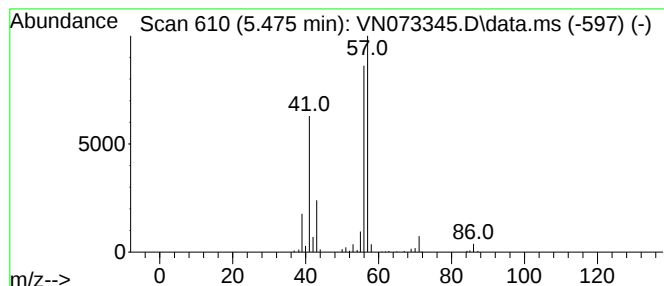
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.475	28.24 ug/l	877572	Pentafluorobenzene	8.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
3			3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	42
4			Butane, 1-chloro-	92	C4H9Cl	000109-69-3	42
5			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	39



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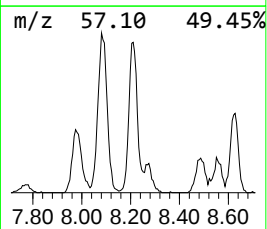
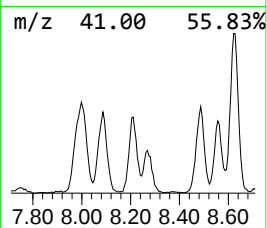
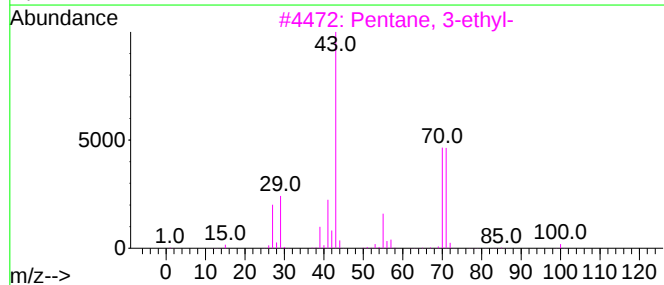
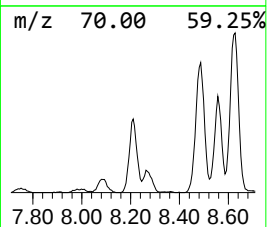
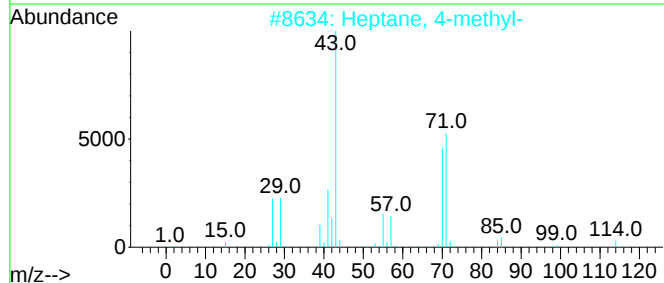
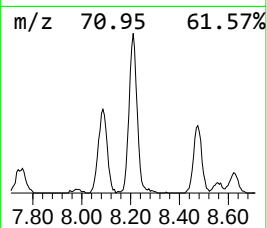
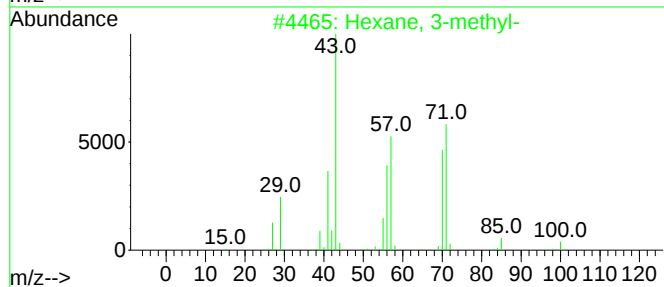
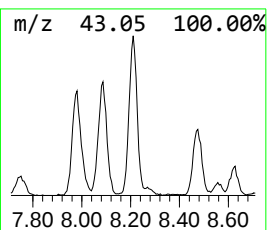
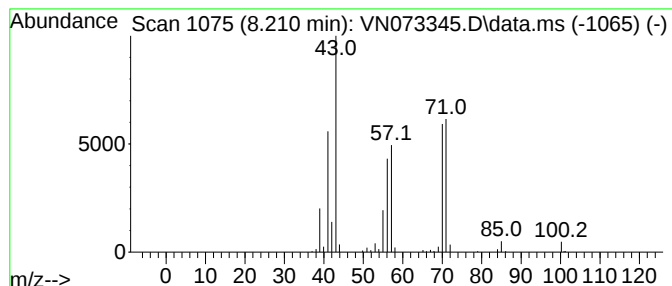
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Hexane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.210	16.94 ug/l	526539	Pentafluorobenzene	8.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	94
2			Heptane, 4-methyl-	114	C8H18	000589-53-7	59
3			Pentane, 3-ethyl-	100	C7H16	000617-78-7	53
4			Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	53
5			Heptane	100	C7H16	000142-82-5	53



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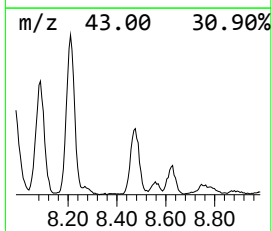
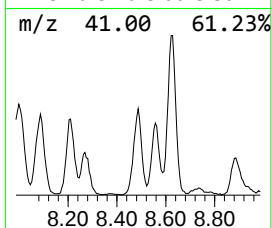
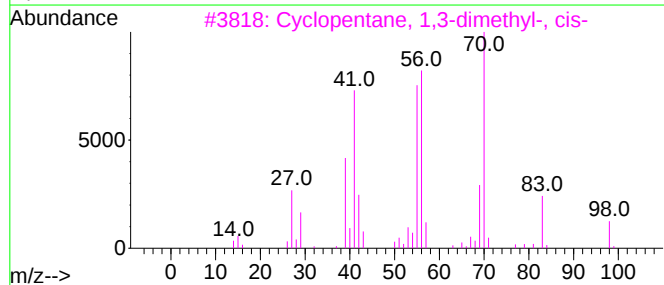
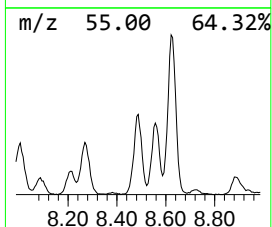
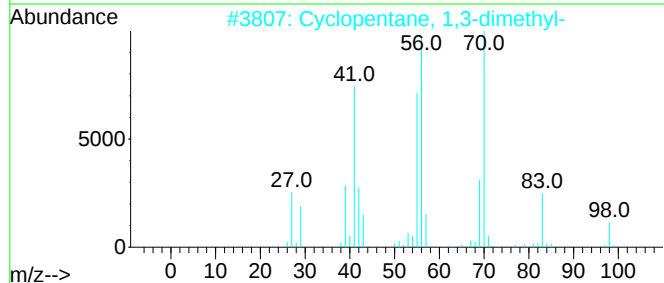
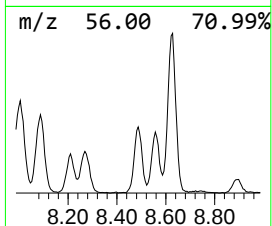
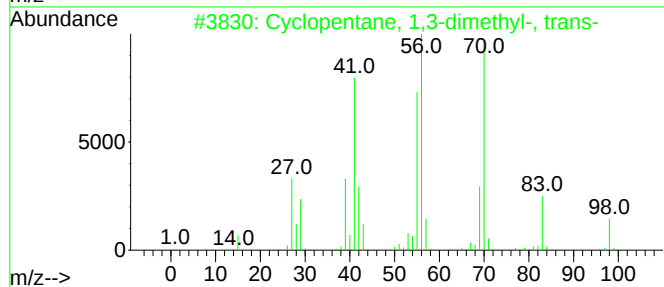
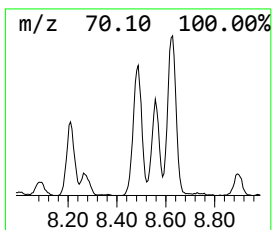
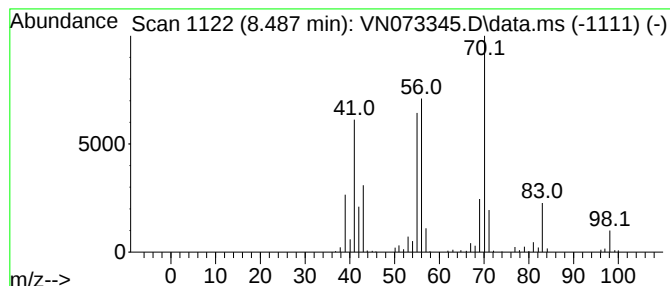
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Quant Title : SW846 8260

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

Peak Number 3 Cyclopentane, 1,3-dimethyl-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.487	22.05 ug/l	633336	1,4-Difluorobenzene	8.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	94
2			Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	94
3			Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	92
4			Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	64
5			Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070622\
Data File : VN073345.D
Acq On : 06 Jul 2022 15:11
Operator : JC\MD
Sample : N3564-04
Misc : 4.97g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-14B-11-1-GR

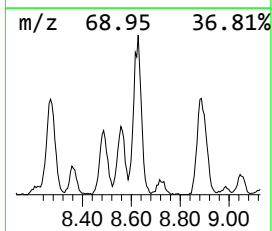
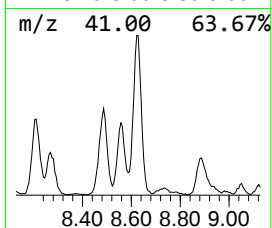
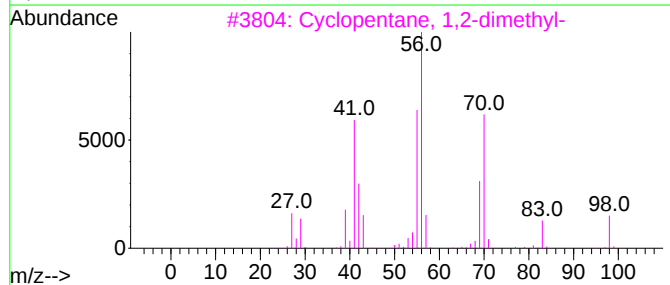
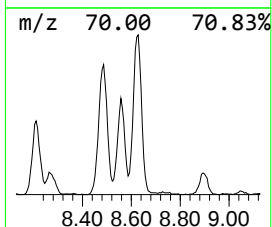
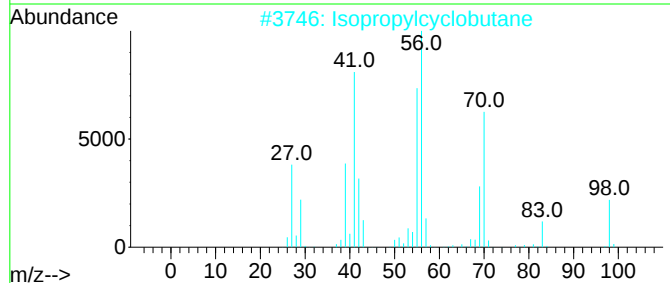
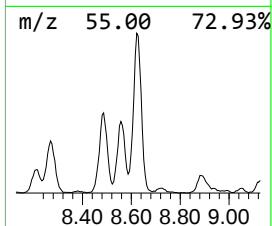
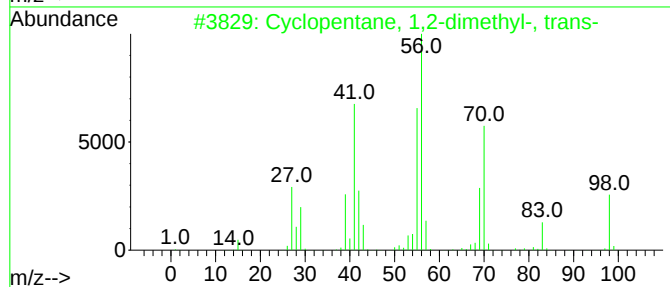
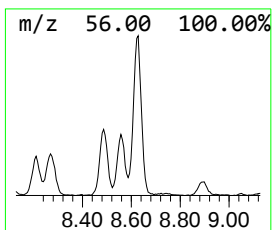
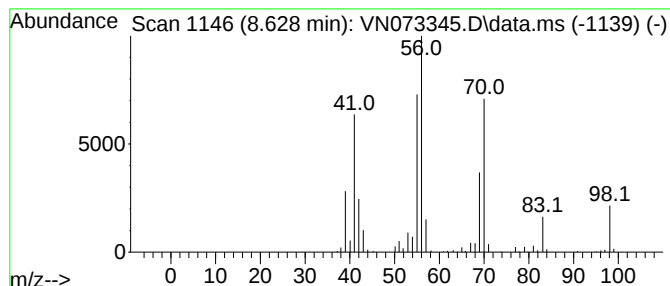
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Quant Title : SW846 8260

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

Peak Number 4 Cyclopentane, 1,2-dimethyl-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.628	36.34 ug/l	1043640	1,4-Difluorobenzene	8.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	94
2			Isopropylcyclobutane	98	C7H14	000872-56-0	94
3			Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	91
4			Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	91
5			1-Heptene	98	C7H14	000592-76-7	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070622\
Data File : VN073345.D
Acq On : 06 Jul 2022 15:11
Operator : JC\MD
Sample : N3564-04
Misc : 4.97g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-14B-11-1-GR

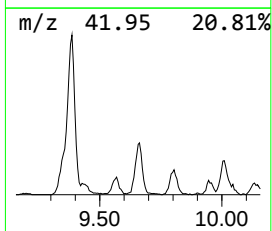
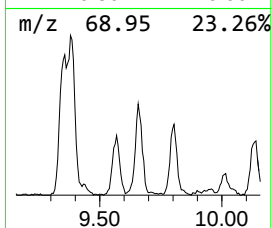
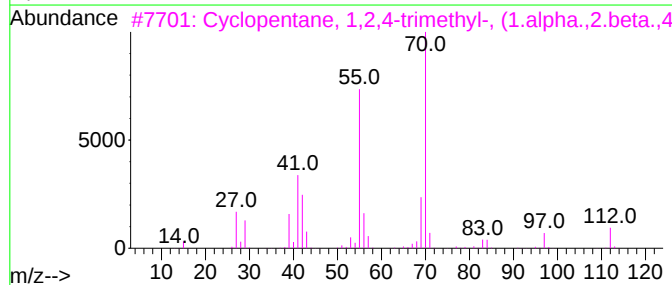
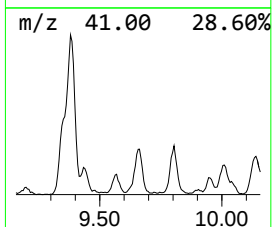
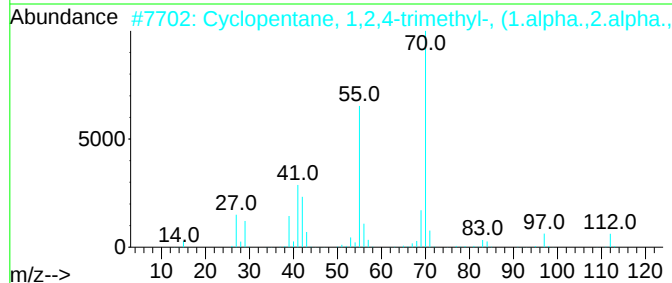
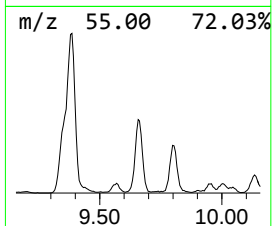
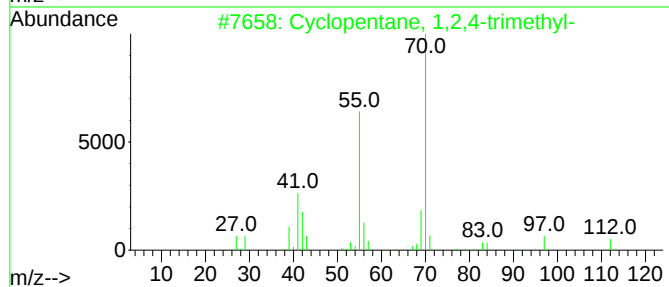
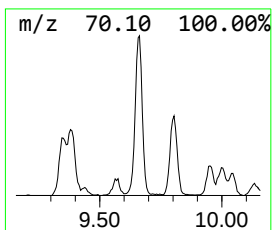
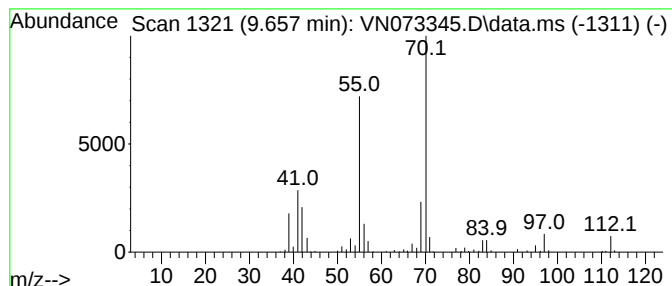
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentane, 1,2,4-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.657	18.75 ug/l	538552	1,4-Difluorobenzene	8.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,2,4-trimethyl-	112	C8H16	002815-58-9	91
2			Cyclopentane, 1,2,4-trimethyl-, ...	112	C8H16	004850-28-6	91
3			Cyclopentane, 1,2,4-trimethyl-, ...	112	C8H16	016883-48-0	87
4			Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	002613-69-6	80
5			Undecane, 3-methylene-	168	C12H24	071138-64-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070622\
Data File : VN073345.D
Acq On : 06 Jul 2022 15:11
Operator : JC\MD
Sample : N3564-04
Misc : 4.97g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-14B-11-1-GR

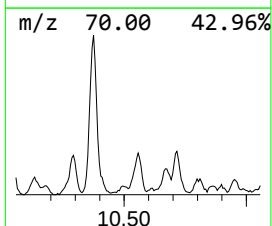
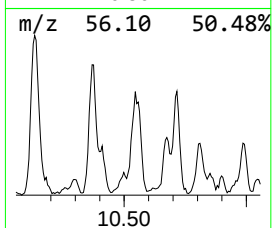
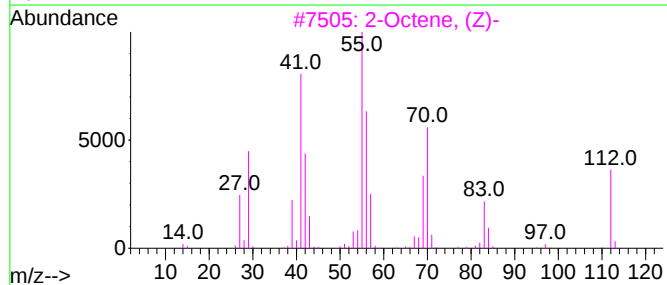
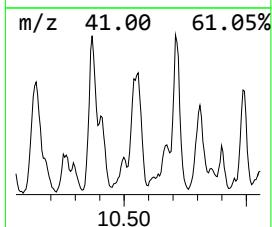
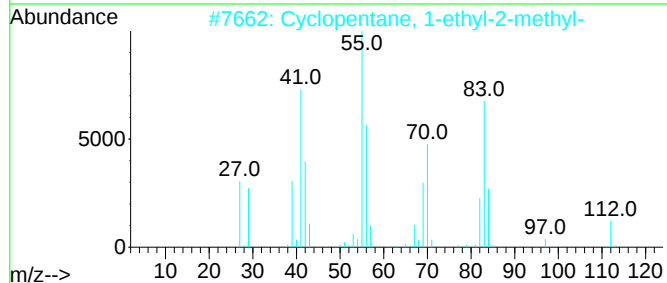
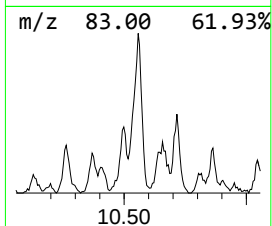
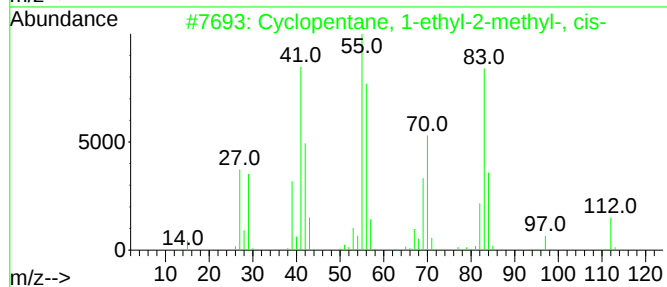
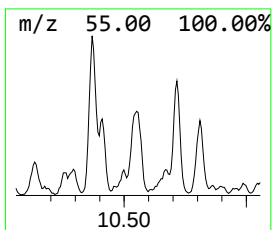
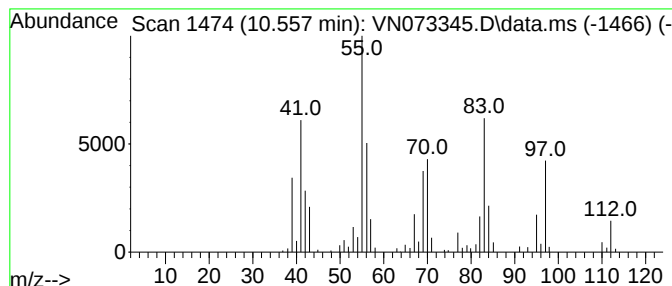
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Quant Title : SW846 8260

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

Peak Number 6 Cyclopentane, 1-ethyl-2-met... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.557	17.60 ug/l	511573	Chlorobenzene-d5	11.681

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1-ethyl-2-methyl-,...	112	C8H16	000930-89-2	94
2			Cyclopentane, 1-ethyl-2-methyl-	112	C8H16	003726-46-3	76
3			2-Octene, (Z)-	112	C8H16	007642-04-8	53
4			4-Octene, (E)-	112	C8H16	014850-23-8	53
5			Cyclooctane	112	C8H16	000292-64-8	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070622\
Data File : VN073345.D
Acq On : 06 Jul 2022 15:11
Operator : JC\MD
Sample : N3564-04
Misc : 4.97g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-14B-11-1-GR

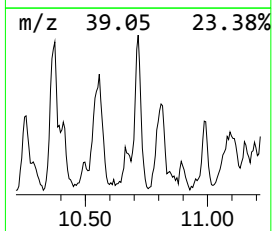
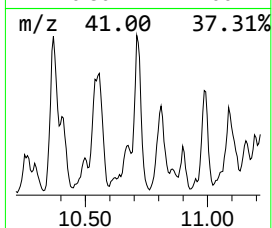
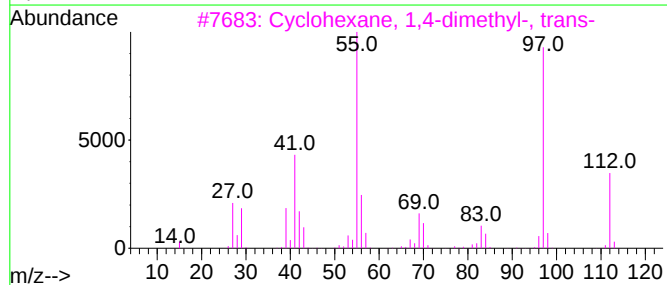
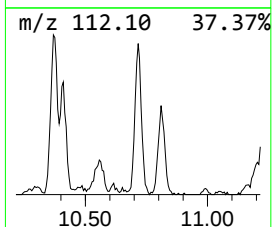
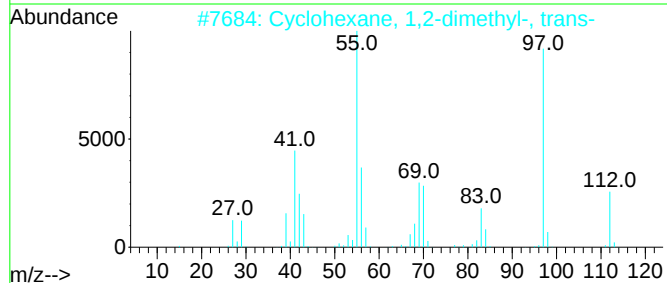
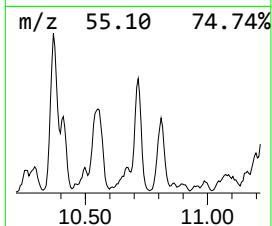
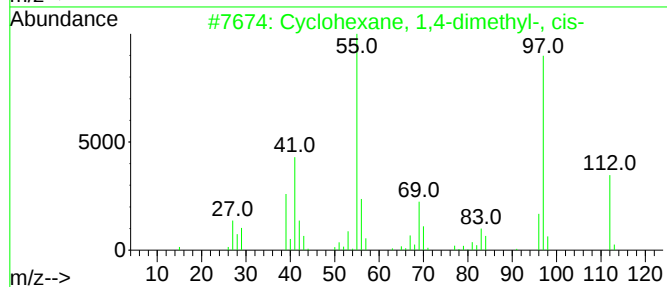
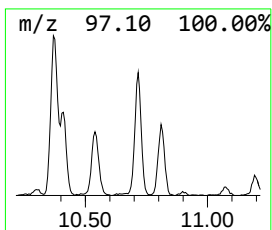
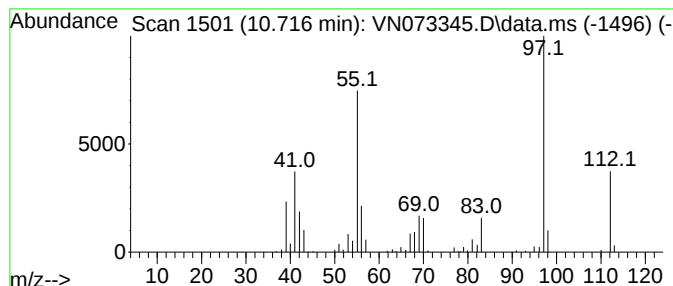
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Quant Title : SW846 8260

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

Peak Number 7 Cyclohexane, 1,4-dimethyl-,... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.716	16.52 ug/l	480302	Chlorobenzene-d5	11.681

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	91
2			Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	91
3			Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	91
4			Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	78
5			1H-Pyrazole, 4,5-dihydro-3,4,5-t...	112	C6H12N2	022591-95-3	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070622\
Data File : VN073345.D
Acq On : 06 Jul 2022 15:11
Operator : JC\MD
Sample : N3564-04
Misc : 4.97g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-14B-11-1-GR

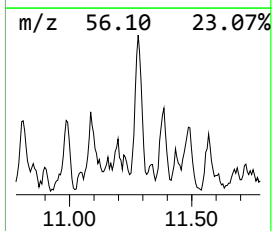
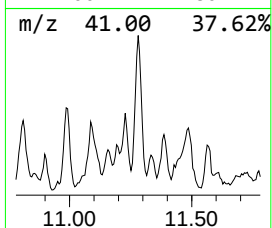
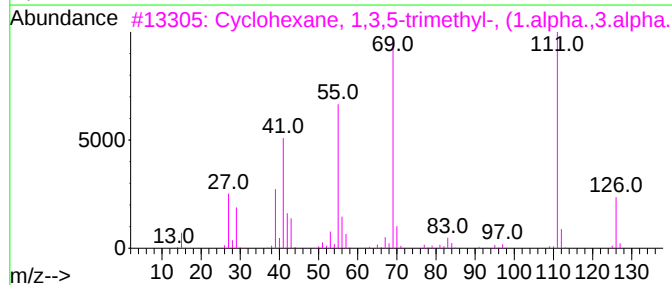
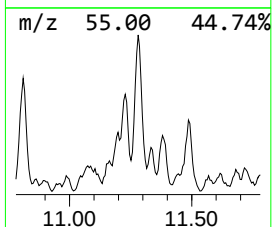
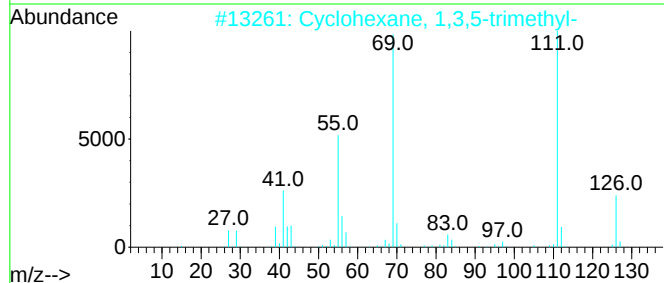
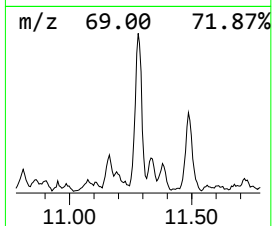
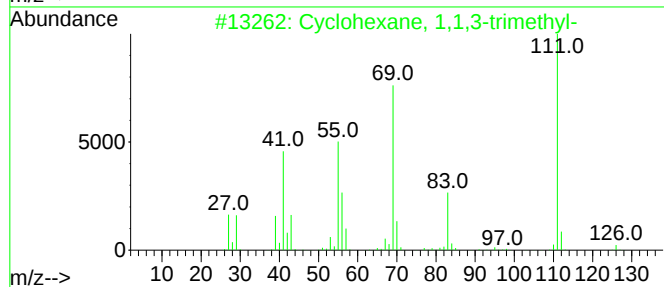
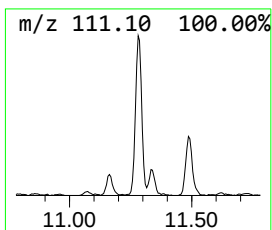
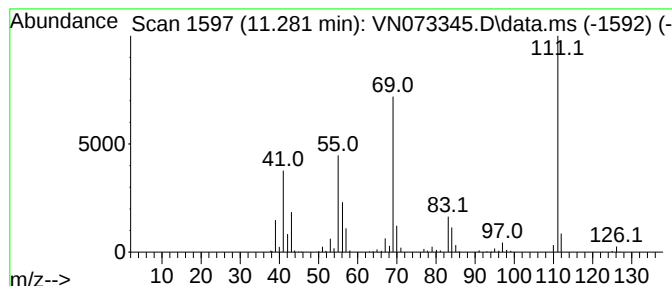
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Quant Title : SW846 8260

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

Peak Number 8 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.281	18.88 ug/l	549005	Chlorobenzene-d5	11.681

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	87
2			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	64
3			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	59
4			4-Amino-6-hydroxypyrimidine	111	C4H5N3O	001193-22-2	53
5			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070622\
Data File : VN073345.D
Acq On : 06 Jul 2022 15:11
Operator : JC\MD
Sample : N3564-04
Misc : 4.97g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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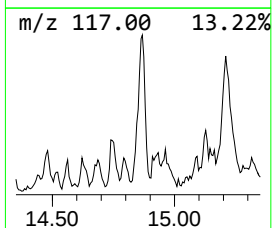
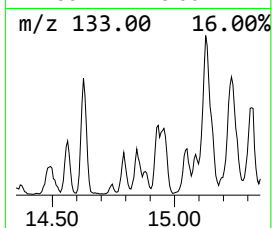
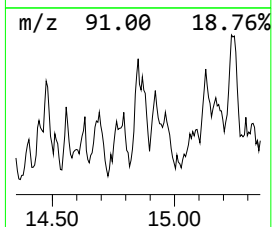
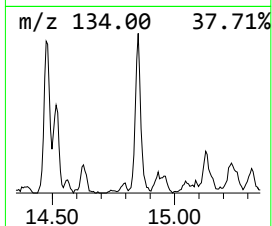
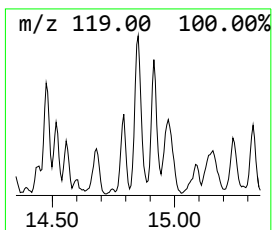
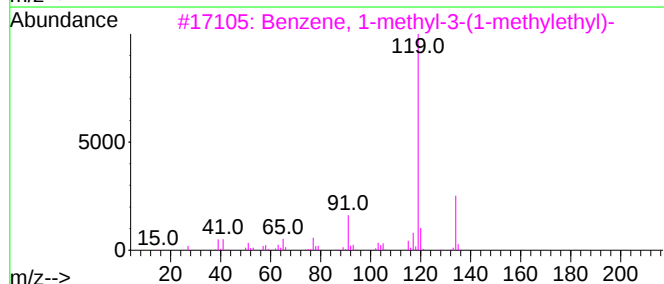
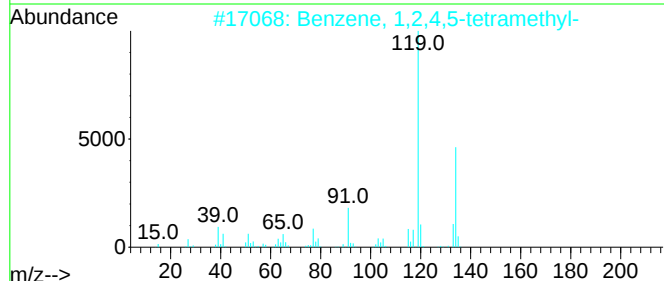
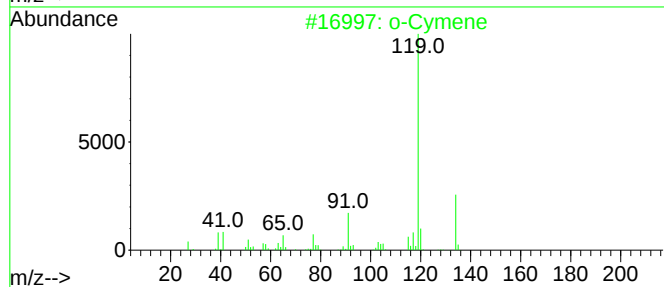
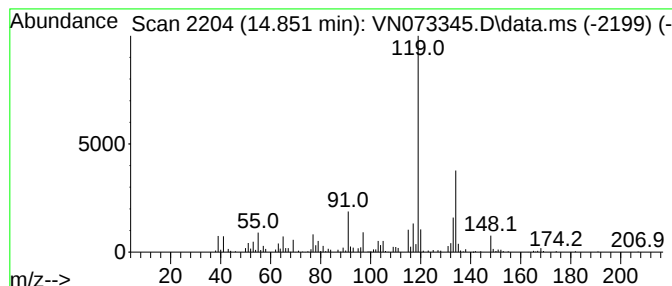
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Quant Title : SW846 8260

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

Peak Number 9 o-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.851	16.51 ug/l	493806	1,4-Dichlorobenzene-d4	13.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	95
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	93
4			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	93
5			p-Cymene	134	C10H14	000099-87-6	89



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070622\
Data File : VN073345.D
Acq On : 06 Jul 2022 15:11
Operator : JC\MD
Sample : N3564-04
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-14B-11-1-GR

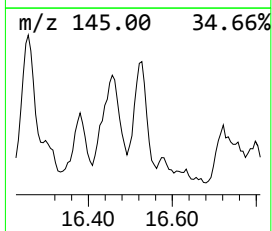
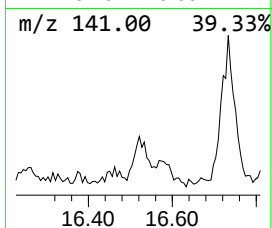
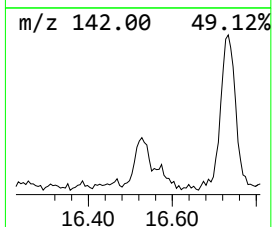
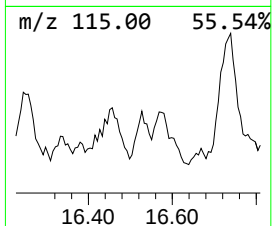
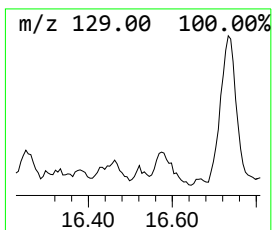
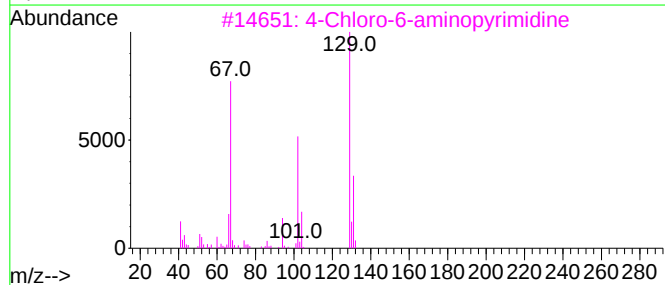
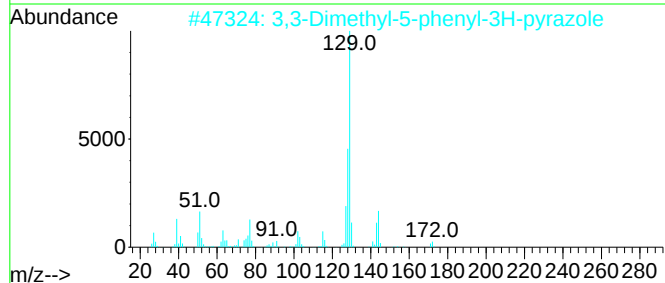
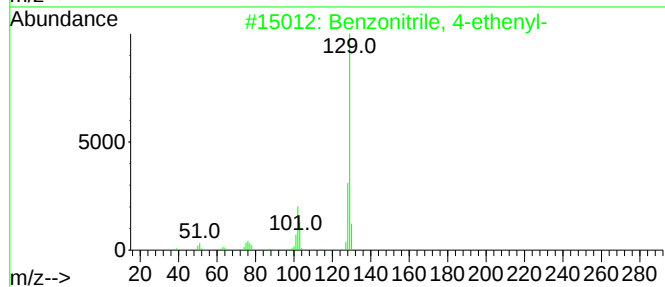
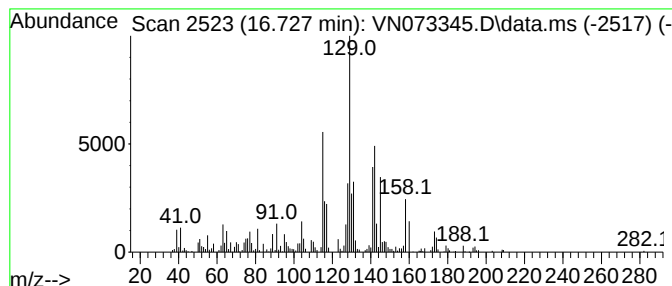
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
Quant Title : SW846 8260

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

Peak Number 10 unknown16.727 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.727	20.82 ug/l	622657	1,4-Dichlorobenzene-d4	13.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzonitrile, 4-ethenyl-	129	C9H7N	003435-51-6	25
2			3,3-Dimethyl-5-phenyl-3H-pyrazole	172	C11H12N2	005363-10-0	25
3			4-Chloro-6-aminopyrimidine	129	C4H4ClN3	005305-59-9	22
4			Cyclohexanecarboxylic acid, 2-(1...	184	C11H20O2	027392-16-1	22
5			2,7-Methanonaphth[2,3-b]oxirene,...	158	C11H10O	013137-34-3	22



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070622\
Data File : VN073345.D
Acq On : 06 Jul 2022 15:11
Operator : JC\MD
Sample : N3564-04
Misc : 4.97g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-14B-11-1-GR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Pentane, 3-methyl-	5.475	28.2	ug/l	877572	1	8.004	1553730	50.0
Hexane, 3-methyl-	8.210	16.9	ug/l	526539	1	8.004	1553730	50.0
Cyclopentane, 1...	8.487	22.1	ug/l	633336	2	8.898	1436100	50.0
Cyclopentane, 1...	8.628	36.3	ug/l	1043640	2	8.898	1436100	50.0
Cyclopentane, 1...	9.657	18.8	ug/l	538552	2	8.898	1436100	50.0
Cyclopentane, 1...	10.557	17.6	ug/l	511573	3	11.681	1453650	50.0
Cyclohexane, 1...	10.716	16.5	ug/l	480302	3	11.681	1453650	50.0
Cyclohexane, 1...	11.281	18.9	ug/l	549005	3	11.681	1453650	50.0
o-Cymene	14.851	16.5	ug/l	493806	4	13.610	1495530	50.0
unknown16.727	16.727	20.8	ug/l	622657	4	13.610	1495530	50.0