

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032024\
 Data File : VN081496.D
 Acq On : 20 Mar 2024 16:10
 Operator : JC\MD
 Sample : P1784-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-5B

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\82N031824W.M

Title : SW846 8260

Signal : TIC: VN081496.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.512	86	95	105	rBV	2592574	4208571	6.51%	0.735%
2	3.248	209	220	240	rBV	13284396	29117024	45.02%	5.087%
3	3.648	278	288	305	rVB2	2062200	5379372	8.32%	0.940%
4	3.859	314	324	339	rVB	764257	1887432	2.92%	0.330%
5	4.048	339	356	364	rBV	410621	1189648	1.84%	0.208%
6	4.148	364	373	391	rVB	1304909	3531148	5.46%	0.617%
7	4.424	404	420	438	rBV2	611446	2140030	3.31%	0.374%
8	5.153	526	544	553	rBV4	253364	987108	1.53%	0.172%
9	5.277	553	565	570	rVV	1956796	6877289	10.63%	1.202%
10	5.353	571	578	608	rVB5	2543159	13888927	21.47%	2.427%
11	5.812	638	656	674	rBV	3380952	11336681	17.53%	1.981%
12	6.312	727	741	754	rBV3	257961	788355	1.22%	0.138%
13	6.653	792	799	812	rVB2	523136	1573062	2.43%	0.275%
14	6.800	812	824	831	rBV2	516430	1616028	2.50%	0.282%
15	6.889	831	839	851	rVV2	414107	1640703	2.54%	0.287%
16	7.083	863	872	887	rVV3	564761	1638430	2.53%	0.286%
17	7.224	887	896	903	rVV2	361856	989064	1.53%	0.173%
18	7.330	903	914	936	rVB	5786944	17115557	26.46%	2.991%
19	8.006	1021	1029	1039	rVB2	746072	1846623	2.86%	0.323%
20	8.165	1047	1056	1060	rBV	423315	1025296	1.59%	0.179%
21	8.259	1060	1072	1078	rVV2	2761218	9260836	14.32%	1.618%
22	8.330	1078	1084	1095	rVB2	1411538	3652495	5.65%	0.638%
23	8.447	1095	1104	1110	rBV	845766	1881512	2.91%	0.329%
24	8.512	1110	1115	1121	rVB	349893	711987	1.10%	0.124%
25	8.583	1121	1127	1135	rBV2	359237	801706	1.24%	0.140%
26	8.724	1142	1151	1157	rBV4	1469774	3896983	6.03%	0.681%
27	8.788	1157	1162	1168	rVV2	1240963	2670615	4.13%	0.467%
28	8.859	1168	1174	1182	rVB	894027	2005008	3.10%	0.350%
29	9.106	1204	1216	1229	rBV2	1535081	4242201	6.56%	0.741%
30	9.335	1248	1255	1260	rBV	330533	714456	1.10%	0.125%
31	9.600	1296	1300	1317	rVB	2100183	4816831	7.45%	0.842%
32	9.777	1323	1330	1337	rVB	521316	981314	1.52%	0.171%
33	10.100	1378	1385	1391	rBV2	1485621	2891874	4.47%	0.505%
34	10.341	1416	1426	1438	rBV3	417108	1212622	1.87%	0.212%
35	10.453	1438	1445	1450	rBV	554714	1049793	1.62%	0.183%
36	10.571	1457	1465	1470	rBV	2055308	4017946	6.21%	0.702%

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Integrator: RTE

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37	10.630	1470	1475	1484	rVB	4424662	8259047	12.77%	1.443%
38	10.759	1489	1497	1507	rVB5	542637	1484293	2.29%	0.259%
39	10.912	1517	1523	1529	rBV	351301	685395	1.06%	0.120%
40	10.982	1529	1535	1546	rVB2	1090192	2630536	4.07%	0.460%
41	11.865	1678	1685	1692	rBV	1532518	2804746	4.34%	0.490%
42	11.965	1692	1702	1712	rVV	21828528	37181948	57.49%	6.497%
43	12.071	1712	1720	1732	rBV	33159630	62938507	97.31%	10.997%
44	12.400	1765	1776	1786	rBV	12314526	20532548	31.75%	3.588%
45	12.694	1817	1826	1833	rVB	2010454	3330639	5.15%	0.582%
46	12.847	1845	1852	1862	rVB	1053587	1951132	3.02%	0.341%
47	13.035	1874	1884	1890	rBV	8313269	13484597	20.85%	2.356%
48	13.100	1890	1895	1897	rVV	4055696	6453271	9.98%	1.128%
49	13.129	1897	1900	1905	rVV	6309654	10410222	16.10%	1.819%
50	13.176	1905	1908	1915	rVB	1670385	2453977	3.79%	0.429%
51	13.335	1928	1935	1945	rBV	9453824	15709590	24.29%	2.745%
52	13.482	1952	1960	1968	rBV2	39870034	64678868	100.00%	11.301%
53	13.612	1974	1982	1988	rVB3	775708	1918351	2.97%	0.335%
54	13.729	1997	2002	2007	rVB	477032	698848	1.08%	0.122%
55	13.823	2007	2018	2025	rBV	8630747	16296935	25.20%	2.847%
56	13.953	2033	2040	2043	rBV	3702578	5763715	8.91%	1.007%
57	13.994	2043	2047	2051	rVV	6319049	11091965	17.15%	1.938%
58	14.035	2051	2054	2063	rVV2	3464432	5716015	8.84%	0.999%
59	14.118	2063	2068	2073	rVB	752943	1185060	1.83%	0.207%
60	14.182	2073	2079	2084	rBV	2407271	3852705	5.96%	0.673%
61	14.247	2084	2090	2093	rVV	5351068	9422035	14.57%	1.646%
62	14.276	2093	2095	2099	rVV	3767862	4584439	7.09%	0.801%
63	14.329	2099	2104	2110	rVV	7798629	12087901	18.69%	2.112%
64	14.400	2110	2116	2119	rVV	1943537	3244706	5.02%	0.567%
65	14.447	2119	2124	2132	rVB2	6056602	10248081	15.84%	1.791%
66	14.576	2140	2146	2152	rVV2	2222930	3559819	5.50%	0.622%
67	14.653	2152	2159	2163	rVV	4988751	8846628	13.68%	1.546%
68	14.694	2163	2166	2171	rVV	7065005	10438644	16.14%	1.824%
69	14.735	2171	2173	2177	rVV	703207	884465	1.37%	0.155%
70	14.882	2192	2198	2201	rBV2	659696	1016669	1.57%	0.178%
71	14.929	2201	2206	2211	rBV	5129222	8255625	12.76%	1.442%
72	14.970	2211	2213	2218	rVB	825848	997547	1.54%	0.174%
73	15.053	2218	2227	2233	rBV2	7599695	16681062	25.79%	2.915%
74	15.117	2233	2238	2248	rVB3	630638	1445269	2.23%	0.253%
75	15.212	2248	2254	2261	rBV2	844300	1717329	2.66%	0.300%
76	15.317	2261	2272	2277	rBV2	2120745	5664635	8.76%	0.990%

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Title : SW846 8260

77	15.441	2285	2293	2304	rVB2	2184701	5583096	8.63%	0.976%
78	15.641	2321	2327	2334	rVB	2403669	4402506	6.81%	0.769%
79	15.753	2340	2346	2351	rBV2	675935	1334816	2.06%	0.233%
80	15.806	2351	2355	2370	rVB	540815	1256060	1.94%	0.219%
81	15.947	2370	2379	2387	rBV	1061994	2130980	3.29%	0.372%
82	16.129	2399	2410	2420	rBV	641959	1966605	3.04%	0.344%
83	16.341	2438	2446	2464	rVB3	455116	1462652	2.26%	0.256%

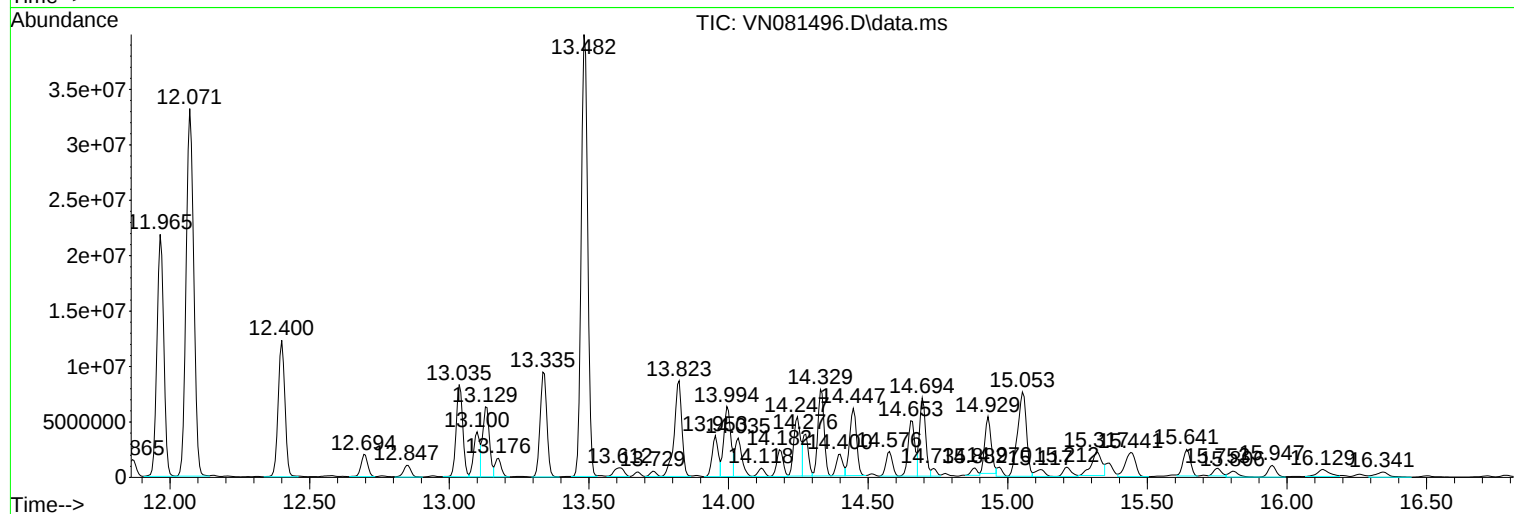
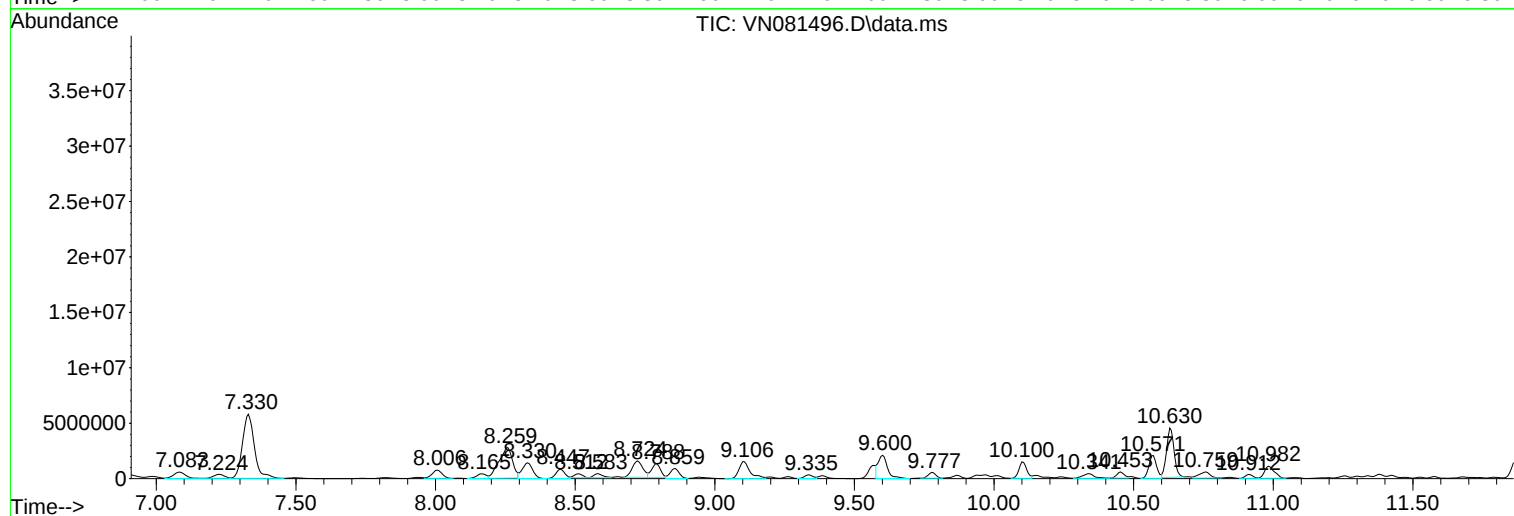
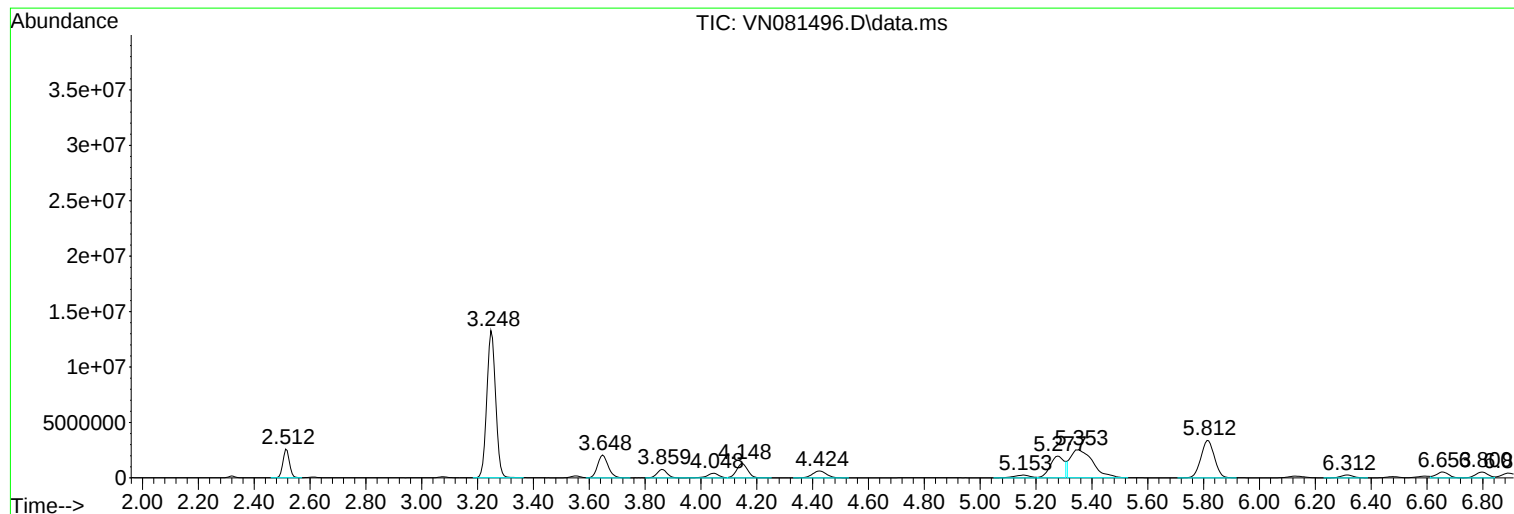
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Instrument :
MSVOA_N
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N031824W.M
Quant Title : SW846 8260

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



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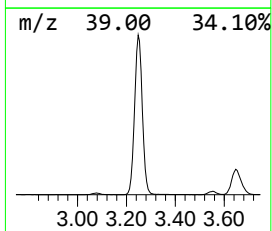
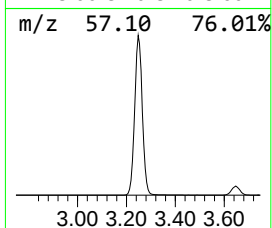
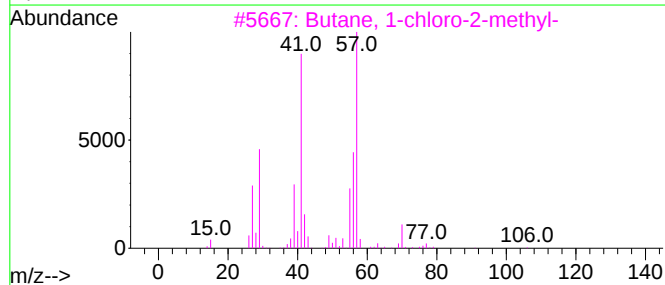
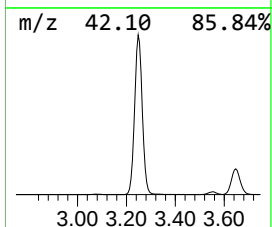
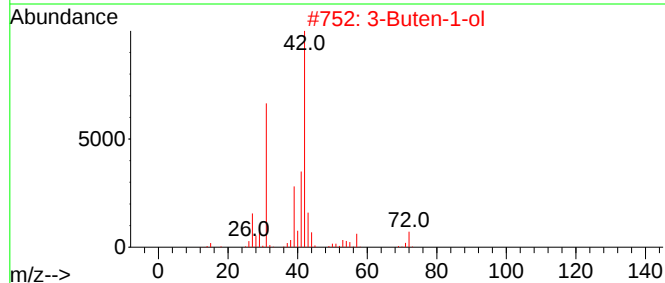
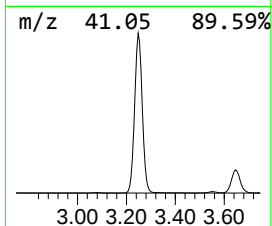
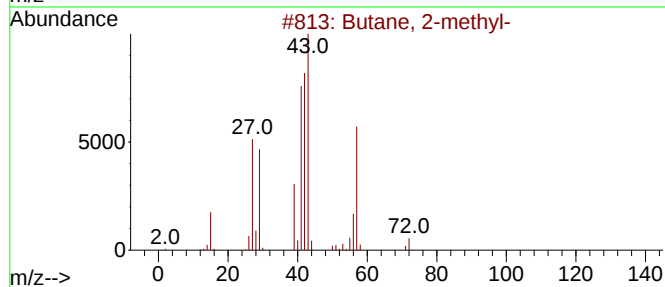
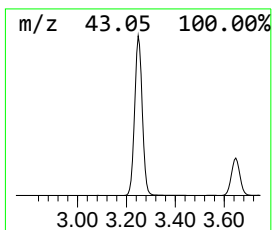
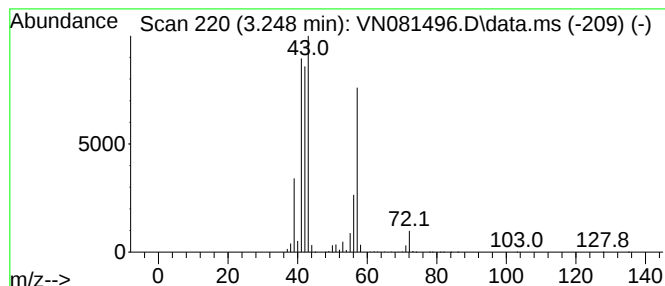
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N031824W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.248	157.21 ug/l	29117000	Pentafluorobenzene	8.224

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-			72	C5H12	000078-78-4	86
2	3-Buten-1-ol			72	C4H8O	000627-27-0	43
3	Butane, 1-chloro-2-methyl-			106	C5H11Cl	000616-13-7	35
4	Pentane			72	C5H12	000109-66-0	28
5	1-Butene			56	C4H8	000106-98-9	10



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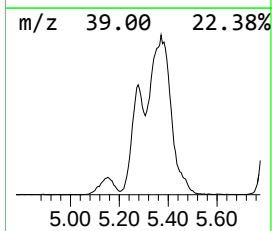
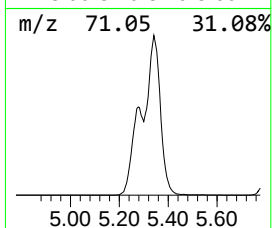
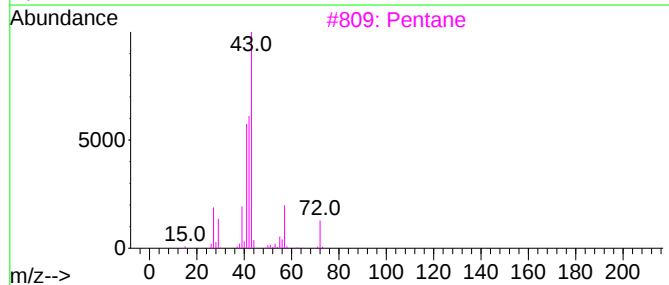
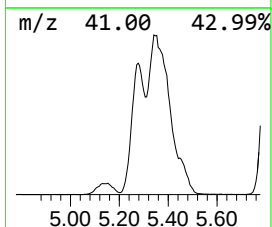
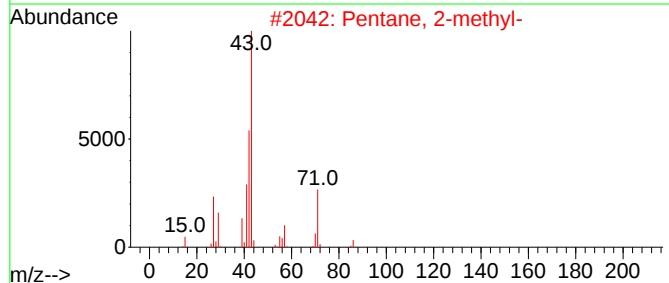
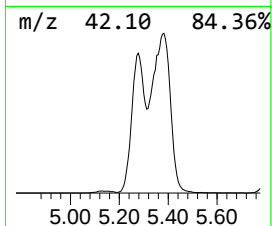
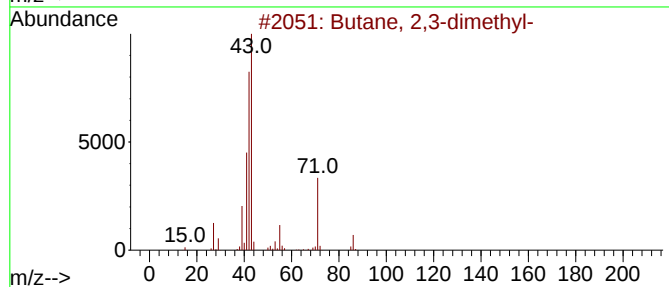
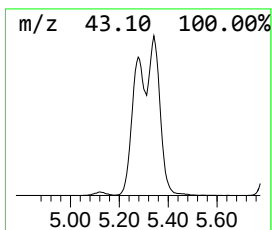
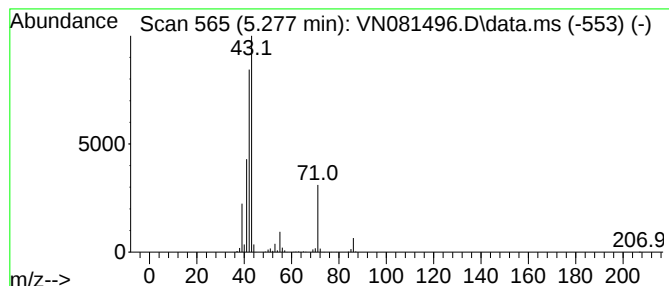
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N031824W.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.277	37.13 ug/l	6877290	Pentafluorobenzene	8.224

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	91
2			Pentane, 2-methyl-	86	C6H14	000107-83-5	50
3			Pentane	72	C5H12	000109-66-0	36
4			1,2,3-Trimethyldiaziridine	86	C4H10N2	113604-56-1	36
5			Pyrrrolidine	71	C4H9N	000123-75-1	27



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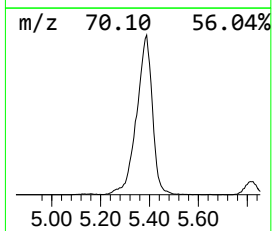
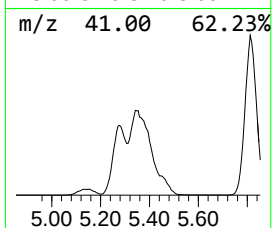
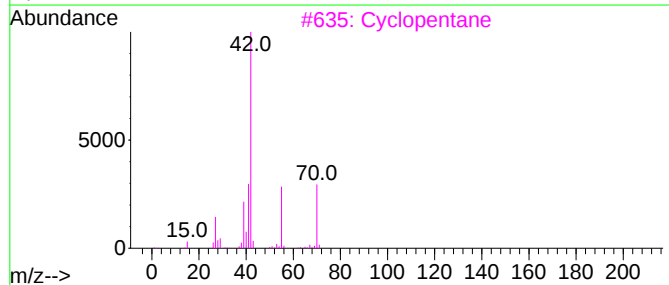
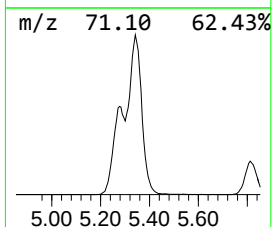
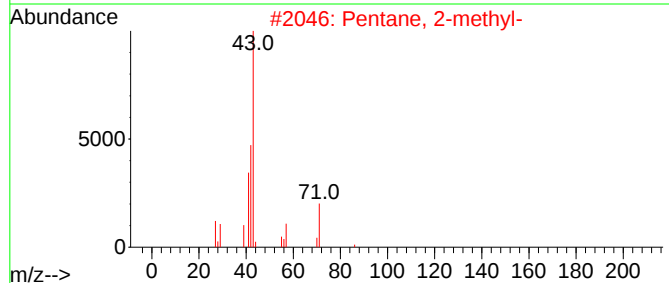
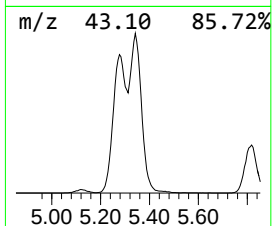
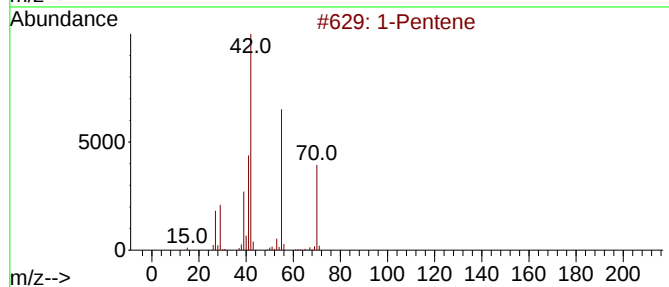
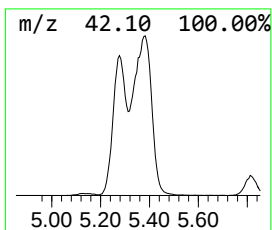
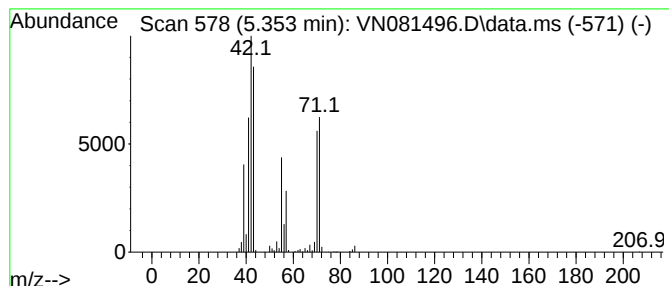
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N031824W.M
Quant Title : SW846 8260

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

Peak Number 3 1-Pentene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.353	74.99 ug/l	13888900	Pentafluorobenzene	8.224

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene			70	C5H10	000109-67-1	50
2	Pentane, 2-methyl-			86	C6H14	000107-83-5	38
3	Cyclopentane			70	C5H10	000287-92-3	37
4	1-Pentanol			88	C5H12O	000071-41-0	32
5	Butane, 2,3-dimethyl-			86	C6H14	000079-29-8	28



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MW-5B

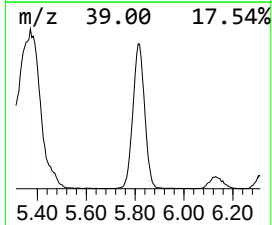
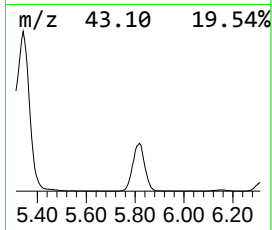
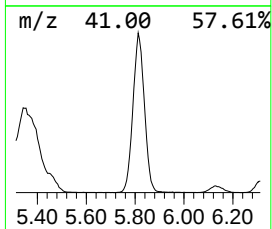
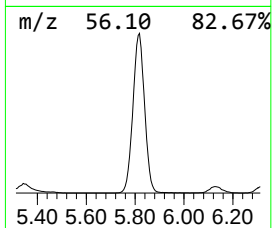
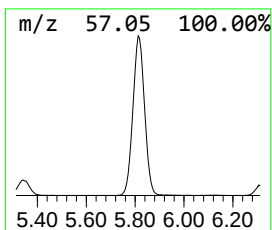
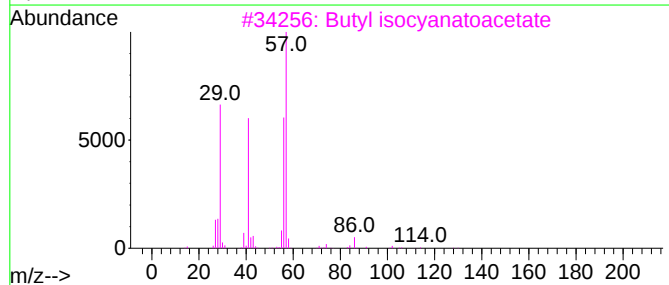
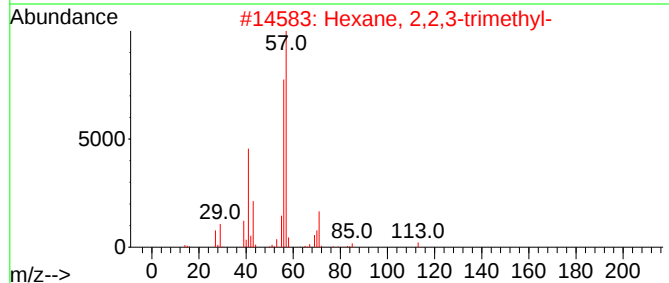
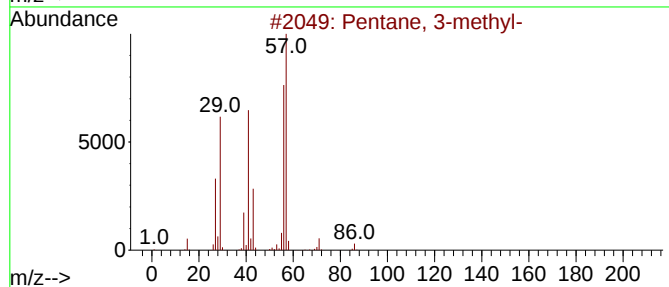
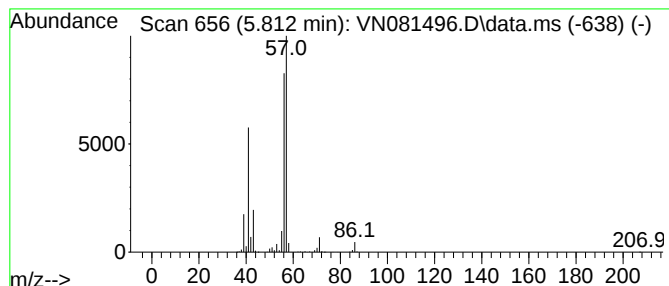
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Quant Title : SW846 8260

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

Peak Number 4 Pentane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.812	61.21 ug/l	11336700	Pentafluorobenzene	8.224

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			Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	64
4			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43
5			Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	40



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032024\
Data File : VN081496.D
Acq On : 20 Mar 2024 16:10
Operator : JC\MD
Sample : P1784-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5B

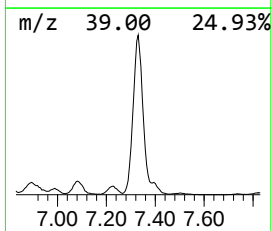
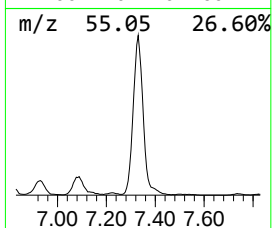
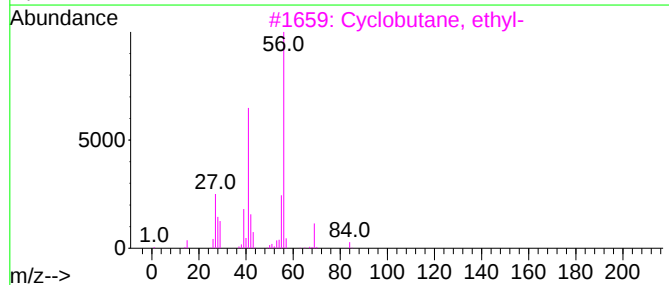
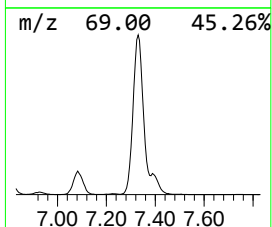
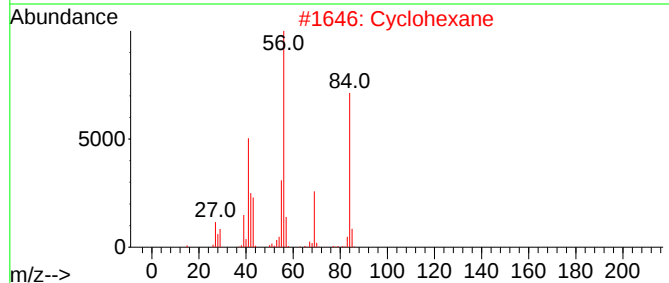
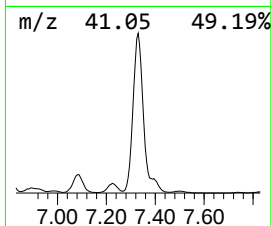
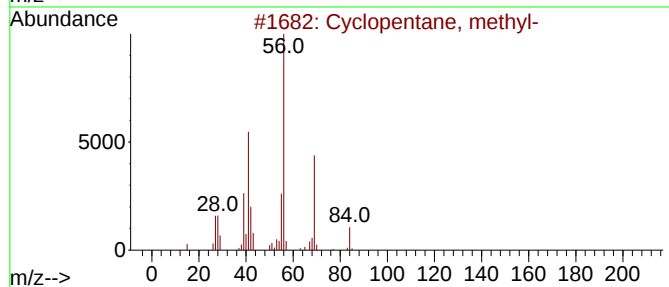
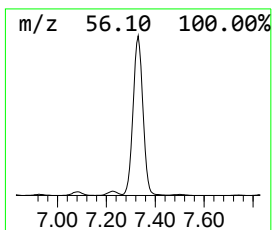
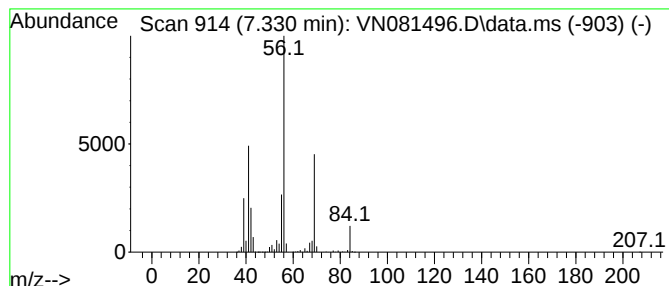
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.330	92.41 ug/l	17115600	Pentafluorobenzene	8.224

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2			Cyclohexane	84	C6H12	000110-82-7	78
3			Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
5			Cyclobutane, butyl-	112	C8H16	013152-44-8	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032024\
Data File : VN081496.D
Acq On : 20 Mar 2024 16:10
Operator : JC\MD
Sample : P1784-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5B

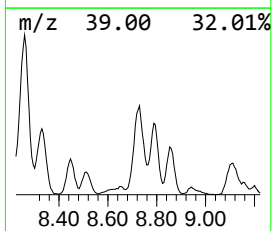
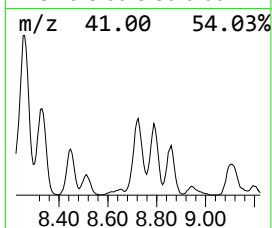
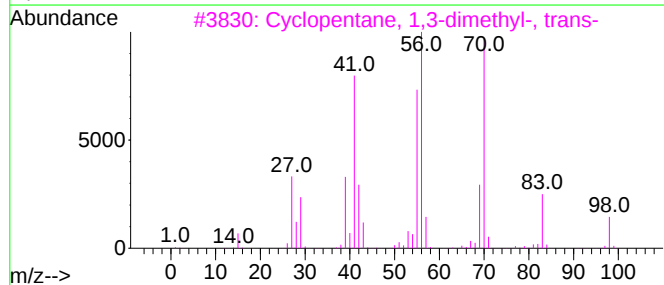
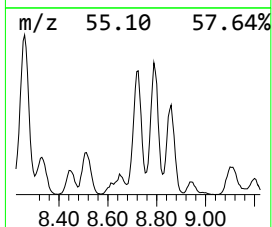
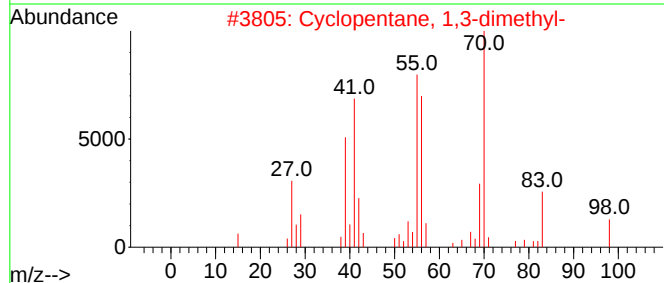
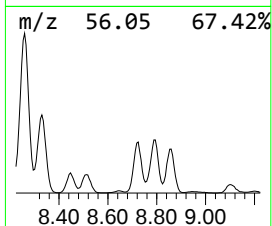
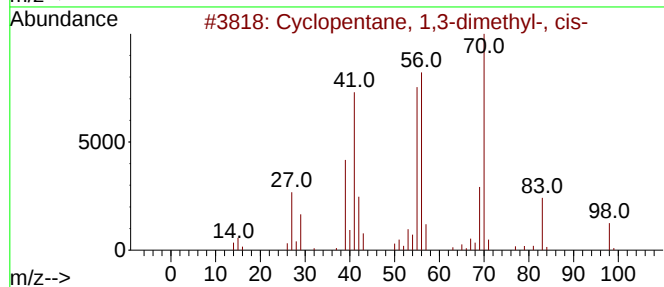
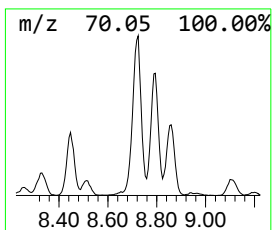
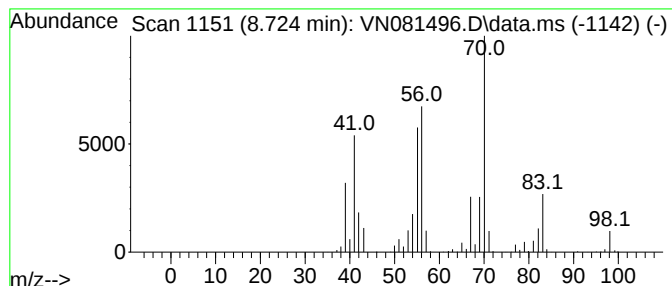
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Quant Title : SW846 8260

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

Peak Number 6 Cyclopentane, 1,3-dimethyl-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.724	45.93 ug/l	3896980	1,4-Difluorobenzene	9.100

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	94
2			Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	80
3			Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	55
4			Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	55
5			Cyclobutane, 2-ethyl-1-methyl-3-...	140	C10H20	061233-72-5	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032024\
Data File : VN081496.D
Acq On : 20 Mar 2024 16:10
Operator : JC\MD
Sample : P1784-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

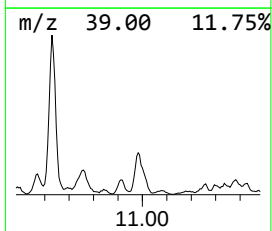
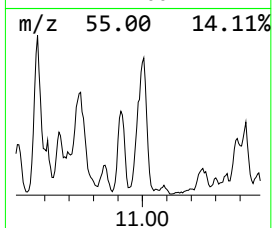
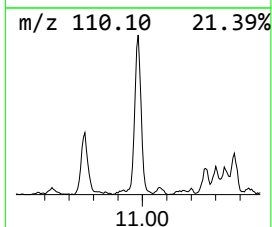
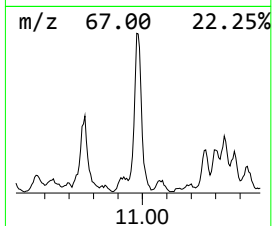
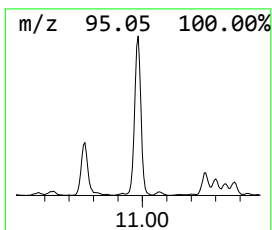
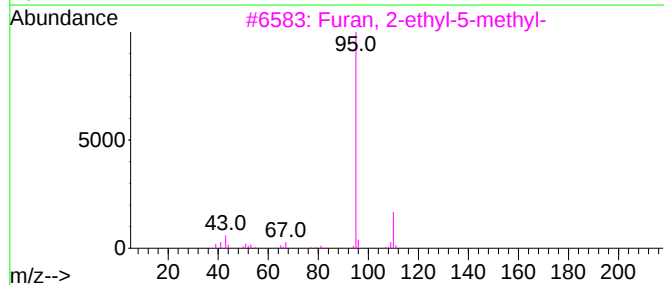
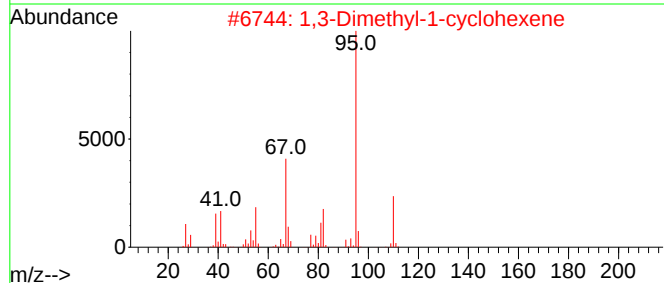
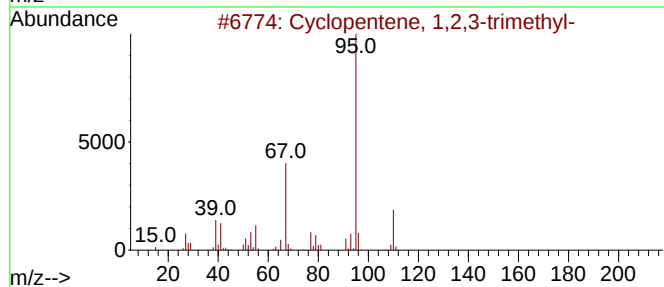
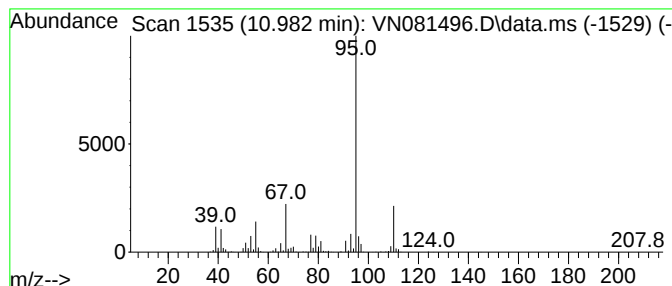
Instrument :
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ClientSampleId :
MW-5B

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N031824W.M
Quant Title : SW846 8260

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

Peak Number 7 Cyclopentene, 1,2,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD			R.T.
10.982	46.89 ug/l	2630540	Chlorobenzene-d5			11.865
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclopentene, 1,2,3-trimethyl-		110	C8H14	000473-91-6	74
2	1,3-Dimethyl-1-cyclohexene		110	C8H14	002808-76-6	72
3	Furan, 2-ethyl-5-methyl-		110	C7H10O	001703-52-2	72
4	2-Isopropylimidazole		110	C6H10N2	036947-68-9	64
5	Bicyclo[3.1.0]hexane, 1,5-dimethyl-		110	C8H14	1010142-17-5	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032024\
Data File : VN081496.D
Acq On : 20 Mar 2024 16:10
Operator : JC\MD
Sample : P1784-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5B

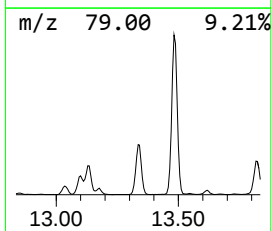
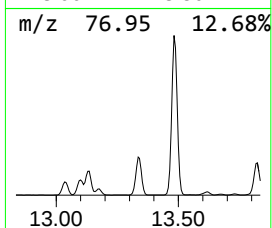
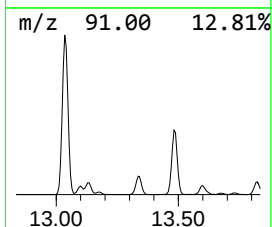
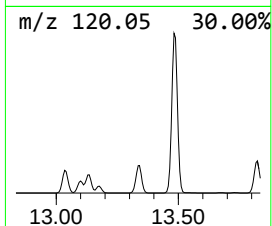
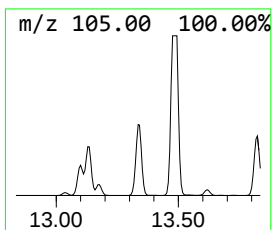
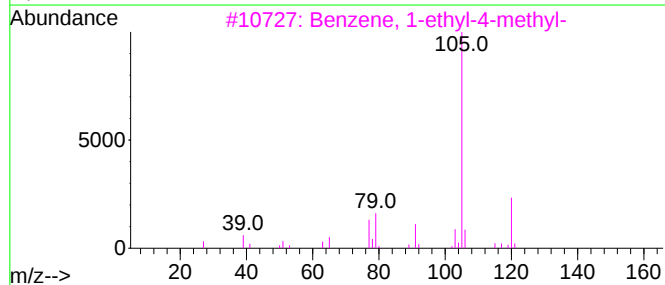
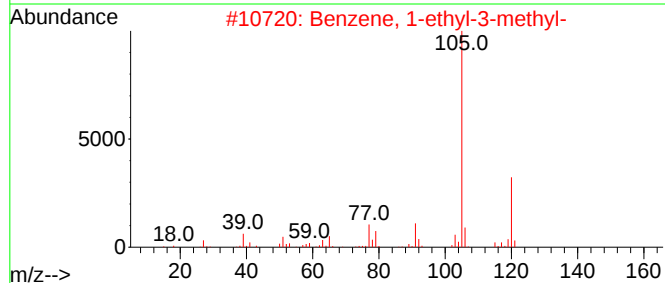
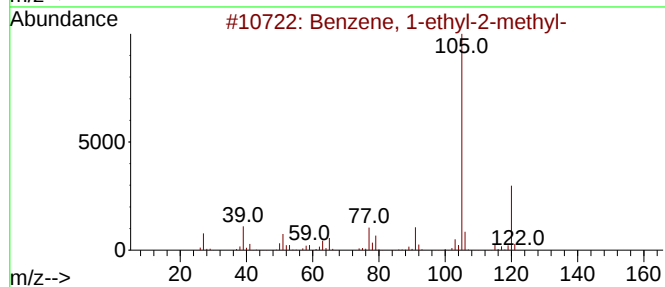
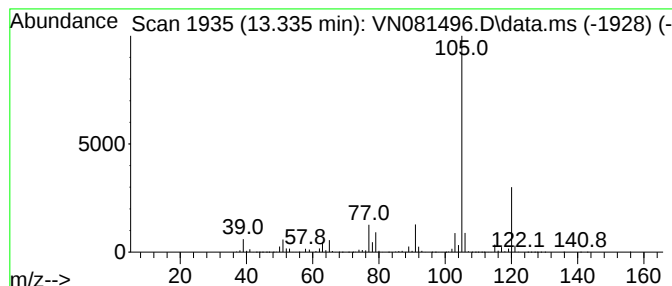
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.335	48.20 ug/l	15709600	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032024\
Data File : VN081496.D
Acq On : 20 Mar 2024 16:10
Operator : JC\MD
Sample : P1784-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5B

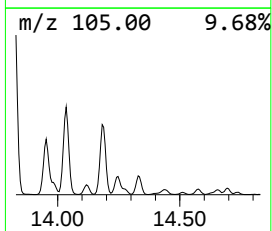
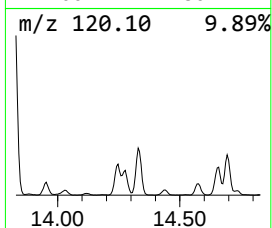
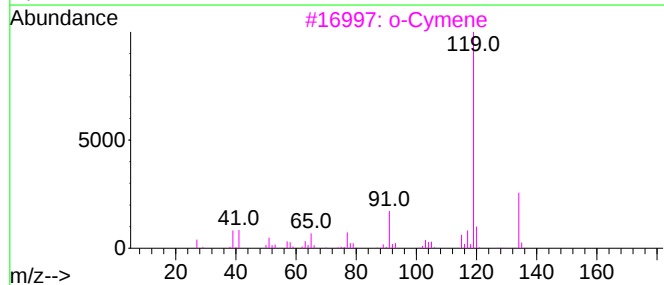
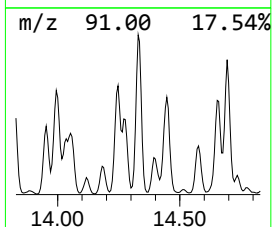
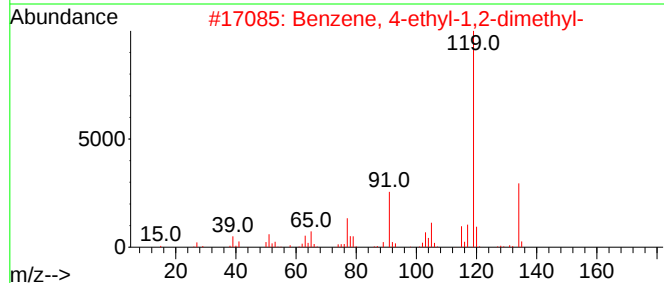
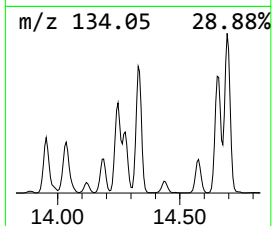
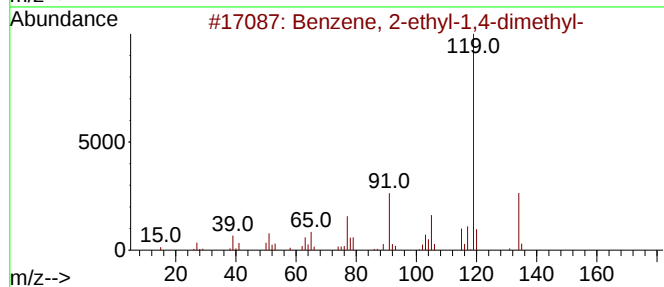
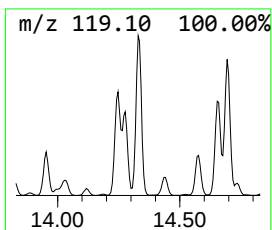
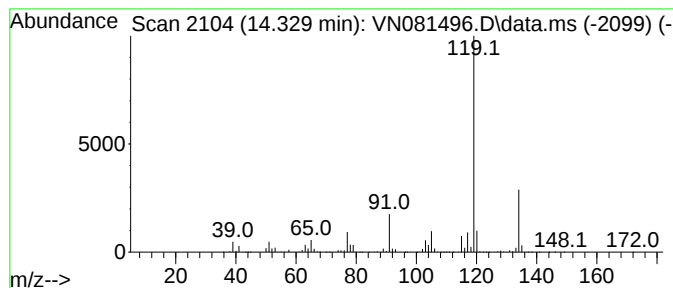
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.329	37.09 ug/l	12087900	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032024\
Data File : VN081496.D
Acq On : 20 Mar 2024 16:10
Operator : JC\MD
Sample : P1784-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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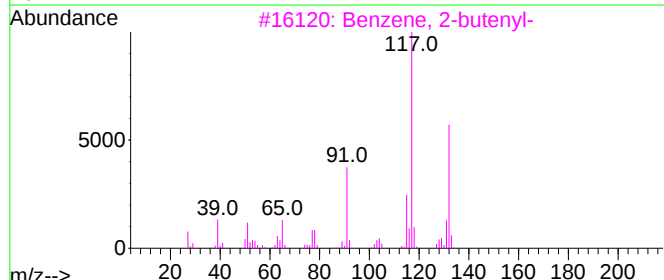
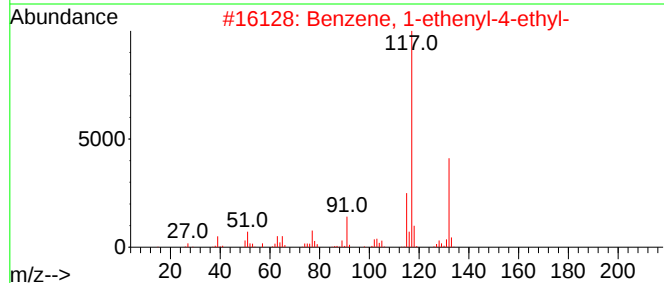
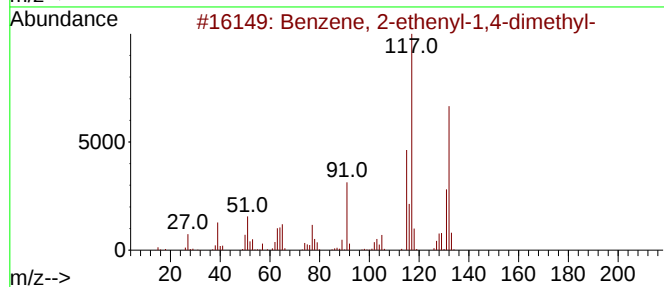
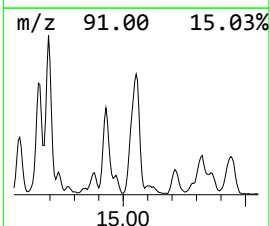
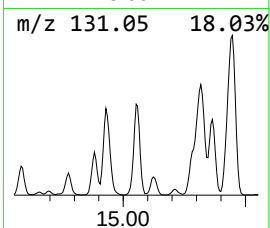
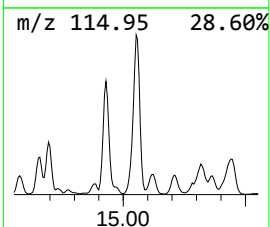
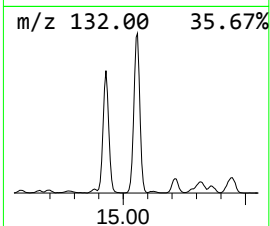
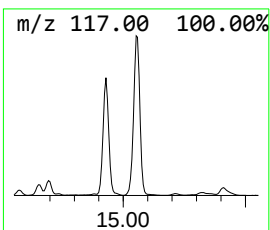
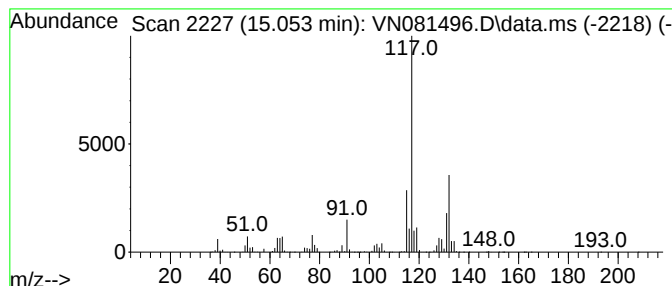
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.053	51.18 ug/l	16681100	1,4-Dichlorobenzene-d4	13.794

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	90
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032024\
Data File : VN081496.D
Acq On : 20 Mar 2024 16:10
Operator : JC\MD
Sample : P1784-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5B

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N031824W.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
Butane, 2-methyl-	3.248	157.2	ug/l	29117000	1	8.224	9260840	50.0
Butane, 2,3-dim...	5.277	37.1	ug/l	6877290	1	8.224	9260840	50.0
1-Pentene	5.353	75.0	ug/l	13888900	1	8.224	9260840	50.0
Pentane, 3-methyl-	5.812	61.2	ug/l	11336700	1	8.224	9260840	50.0
Cyclopentane, m...	7.330	92.4	ug/l	17115600	1	8.224	9260840	50.0
Cyclopentane, 1...	8.724	45.9	ug/l	3896980	2	9.100	4242200	50.0
Cyclopentene, 1...	10.982	46.9	ug/l	2630540	3	11.865	2804750	50.0
Benzene, 1-ethy...	13.335	48.2	ug/l	15709600	4	13.794	16296900	50.0
Benzene, 2-ethy...	14.329	37.1	ug/l	12087900	4	13.794	16296900	50.0
Benzene, 1-ethe...	15.053	51.2	ug/l	16681100	4	13.794	16296900	50.0