

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
 Data File : VN083966.D
 Acq On : 18 Sep 2024 06:45
 Operator : JC\MD
 Sample : P3964-05
 Misc : 10.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SINK-2

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\82N091024W.M
 Title : SW846 8260

Signal : TIC: VN083966.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.253	547	561	564	rBV3	51350	168470	1.22%	0.182%
2	5.312	566	571	587	rVB3	86321	283042	2.05%	0.307%
3	5.794	639	653	666	rVB2	59942	201528	1.46%	0.218%
4	7.218	884	895	903	rBV2	112092	326525	2.36%	0.354%
5	7.318	903	912	920	rBV2	76337	202419	1.46%	0.219%
6	7.994	1021	1027	1035	rVB2	61795	149844	1.08%	0.162%
7	8.165	1046	1056	1058	rBV	132630	284320	2.06%	0.308%
8	8.218	1058	1065	1075	rVV2	512458	1318695	9.54%	1.428%
9	8.324	1075	1083	1093	rVB	221819	574260	4.15%	0.622%
10	8.441	1093	1103	1112	rBV	316516	711619	5.15%	0.771%
11	8.577	1118	1126	1140	rBV	154576	360488	2.61%	0.390%
12	8.753	1140	1156	1168	rBV	1629426	4219603	30.52%	4.569%
13	8.853	1168	1173	1181	rVB4	60273	146333	1.06%	0.158%
14	8.965	1181	1192	1206	rVB2	222587	563409	4.08%	0.610%
15	9.094	1206	1214	1221	rBV3	516886	1195437	8.65%	1.295%
16	9.335	1248	1255	1260	rVV3	92690	234820	1.70%	0.254%
17	9.588	1283	1298	1303	rBV3	621589	1503910	10.88%	1.629%
18	9.641	1303	1307	1314	rVV	387075	715673	5.18%	0.775%
19	9.724	1314	1321	1326	rVV	133543	287275	2.08%	0.311%
20	9.777	1326	1330	1335	rVV2	86790	174479	1.26%	0.189%
21	9.865	1335	1345	1351	rVB5	63548	175522	1.27%	0.190%
22	10.012	1362	1370	1379	rVB	1232217	2430097	17.58%	2.632%
23	10.141	1379	1392	1398	rBV	1708931	3615514	26.15%	3.915%
24	10.200	1398	1402	1406	rBV	202565	293669	2.12%	0.318%
25	10.335	1416	1425	1438	rBV2	336500	863777	6.25%	0.935%
26	10.488	1438	1451	1458	rBV	1097645	2057249	14.88%	2.228%
27	10.565	1458	1464	1469	rVV	631626	1156580	8.37%	1.252%
28	10.629	1469	1475	1484	rVV	313742	678323	4.91%	0.735%
29	10.741	1488	1494	1501	rVB2	338223	647504	4.68%	0.701%
30	10.847	1508	1512	1518	rVB3	80537	146956	1.06%	0.159%
31	10.924	1518	1525	1530	rBV4	61429	152582	1.10%	0.165%
32	11.006	1530	1539	1546	rVV4	219119	531382	3.84%	0.575%
33	11.082	1546	1552	1558	rVB5	76531	150411	1.09%	0.163%
34	11.171	1559	1567	1573	rBV2	93813	198933	1.44%	0.215%
35	11.271	1574	1584	1592	rBV	150418	348675	2.52%	0.378%
36	11.641	1639	1647	1659	rVB3	302655	727537	5.26%	0.788%

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Integrator: RTE

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Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
 Title : SW846 8260

37	11.741	1659	1664	1674	rBV	233068	403970	2.92%	0.437%
38	11.865	1675	1685	1691	rBV2	574950	1280512	9.26%	1.387%
39	11.965	1691	1702	1710	rVB	1662919	3053860	22.09%	3.307%
40	12.065	1711	1719	1731	rBV	7148049	13824328	100.00%	14.971%
41	12.394	1766	1775	1784	rVV	3049723	5372059	38.86%	5.818%
42	12.476	1784	1789	1797	rVB	95703	189343	1.37%	0.205%
43	12.694	1820	1826	1832	rBV	646955	1138101	8.23%	1.232%
44	12.823	1841	1848	1850	rBV	339619	608685	4.40%	0.659%
45	12.988	1871	1876	1878	rBV2	115077	184072	1.33%	0.199%
46	13.035	1878	1884	1889	rVV	1016145	1712466	12.39%	1.854%
47	13.094	1889	1894	1898	rVV	3413082	5851086	42.32%	6.336%
48	13.129	1898	1900	1903	rVV	1497832	1961072	14.19%	2.124%
49	13.170	1903	1907	1918	rVB	1743166	3006271	21.75%	3.256%
50	13.335	1927	1935	1942	rBV	1175608	1993165	14.42%	2.158%
51	13.423	1942	1950	1954	rBV	244707	460251	3.33%	0.498%
52	13.482	1954	1960	1966	rBV	5084395	8076019	58.42%	8.746%
53	13.529	1966	1968	1973	rVB	122349	159997	1.16%	0.173%
54	13.641	1973	1987	1991	rBV4	357486	1112569	8.05%	1.205%
55	13.817	2007	2017	2033	rVB3	1292791	3977642	28.77%	4.307%
56	13.988	2040	2046	2049	rBV2	869818	1703980	12.33%	1.845%
57	14.106	2062	2066	2074	rVB3	140905	246036	1.78%	0.266%
58	14.182	2074	2079	2084	rVB	238790	410256	2.97%	0.444%
59	14.241	2084	2089	2092	rBV	396041	665943	4.82%	0.721%
60	14.270	2092	2094	2099	rVV	354332	440164	3.18%	0.477%
61	14.329	2099	2104	2112	rVB	663369	1052022	7.61%	1.139%
62	14.441	2118	2123	2128	rVB3	180123	297096	2.15%	0.322%
63	14.570	2141	2145	2151	rVB2	144259	240427	1.74%	0.260%
64	14.653	2151	2159	2162	rBV	412572	705750	5.11%	0.764%
65	14.688	2162	2165	2169	rVV	486354	696352	5.04%	0.754%
66	14.735	2169	2173	2177	rVB2	161237	228277	1.65%	0.247%
67	14.923	2200	2205	2209	rBV	227482	421289	3.05%	0.456%
68	15.047	2216	2226	2232	rBV4	340862	1125190	8.14%	1.218%
69	15.312	2267	2271	2283	rVB2	146069	372888	2.70%	0.404%
70	15.429	2286	2291	2300	rVB4	93560	241462	1.75%	0.261%
71	15.635	2314	2326	2338	rVB	353394	854435	6.18%	0.925%
72	15.941	2372	2378	2385	rBV2	71940	158016	1.14%	0.171%
73	16.776	2513	2520	2521	rBV	206693	321163	2.32%	0.348%

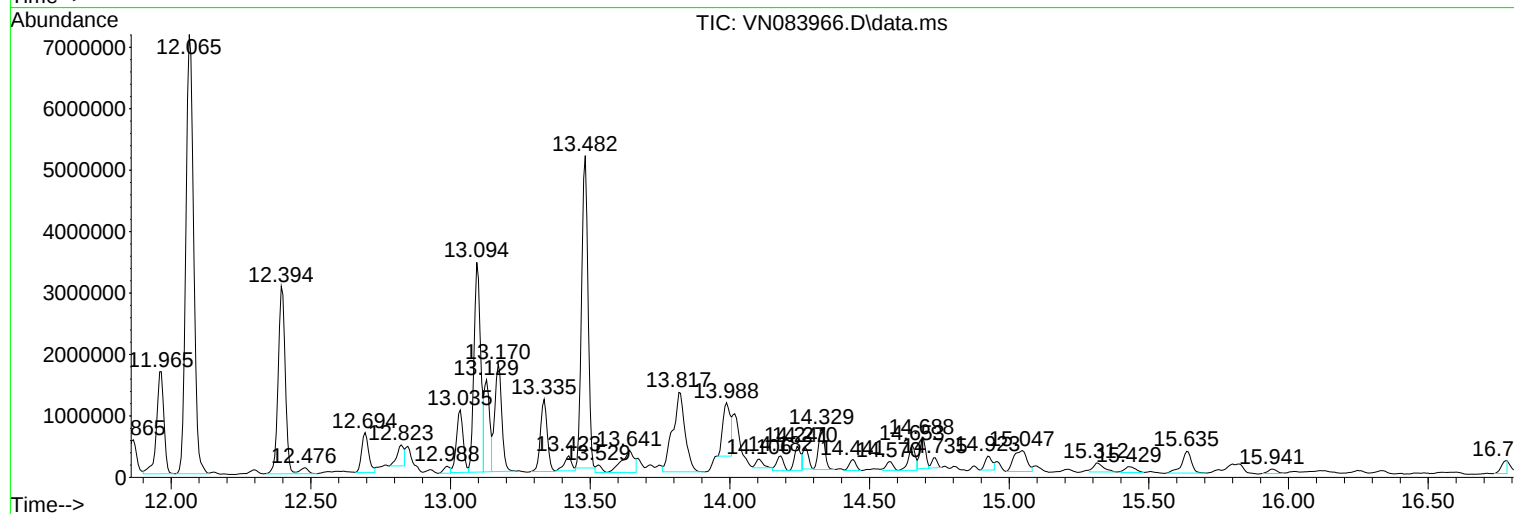
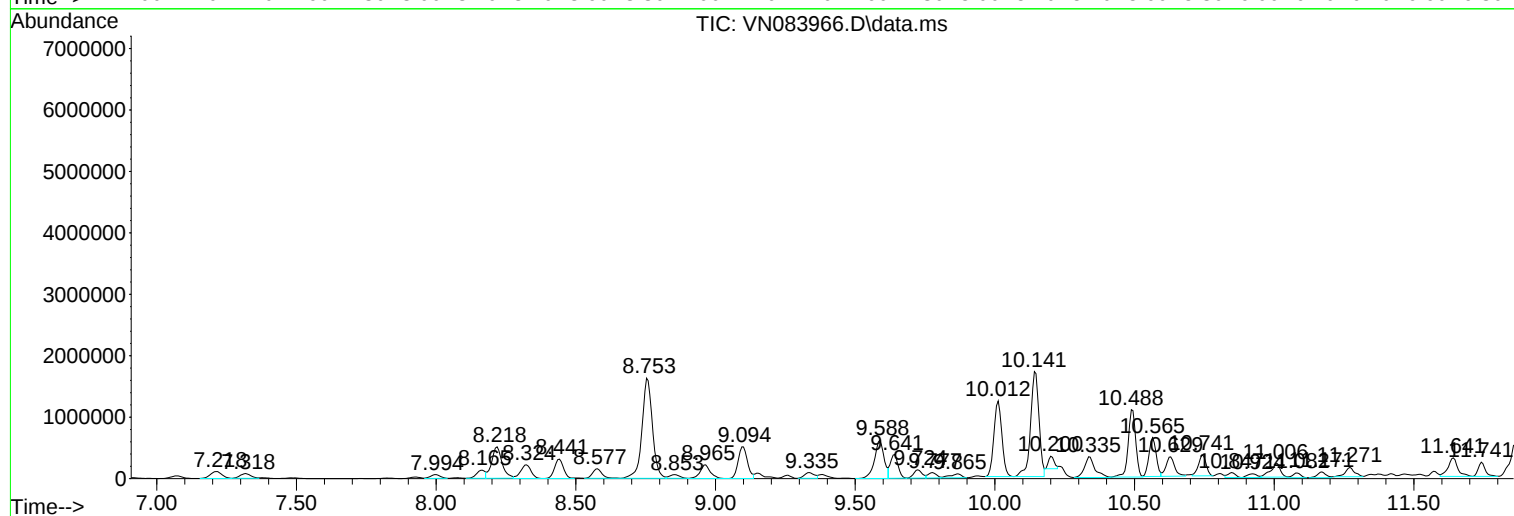
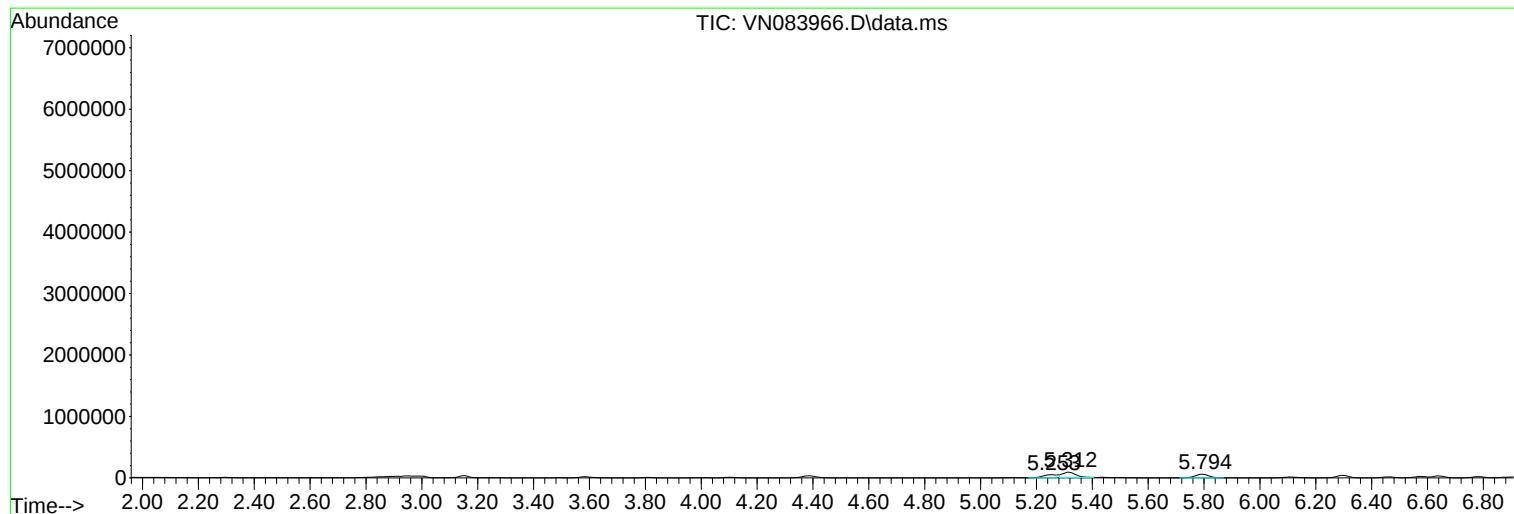
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ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

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



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Sample : P3964-05
Misc : 10.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-2

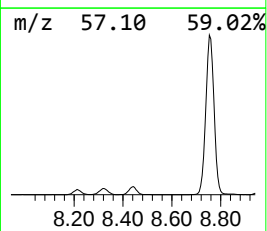
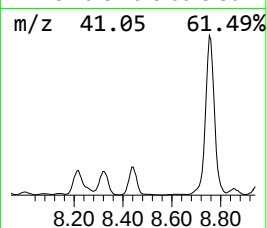
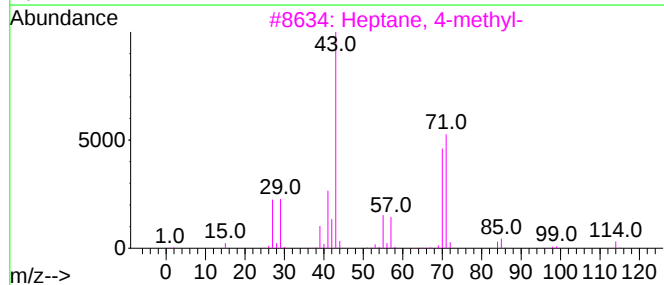
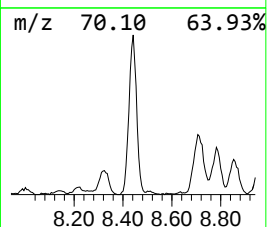
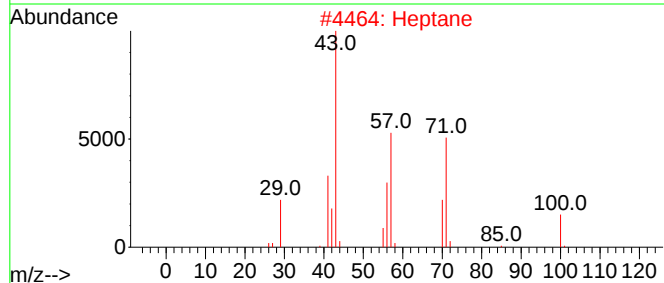
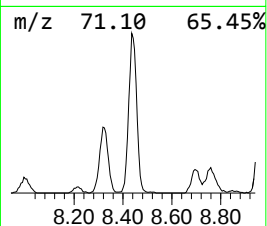
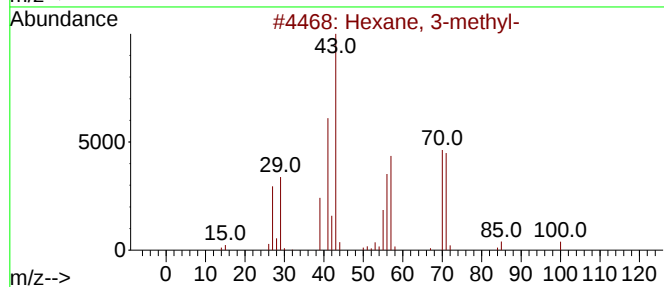
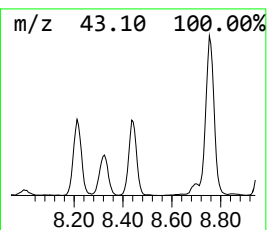
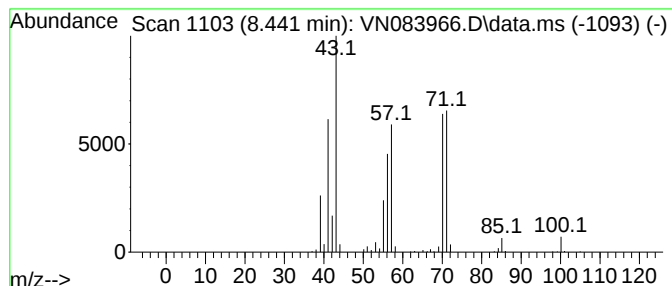
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Quant Title : SW846 8260

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

Peak Number 1 Hexane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.441	26.98 ug/l	711619	Pentafluorobenzene	8.218

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	95
2			Heptane	100	C7H16	000142-82-5	80
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53
5			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	53



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Misc : 10.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-2

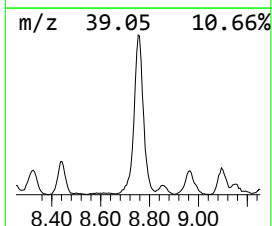
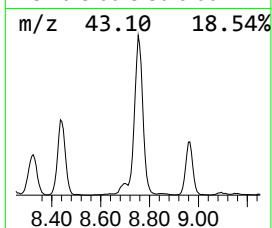
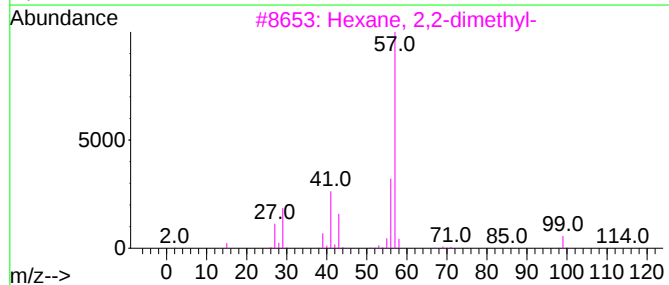
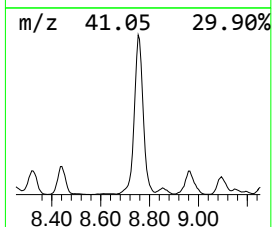
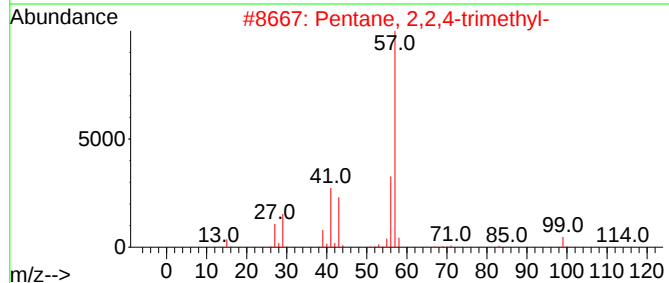
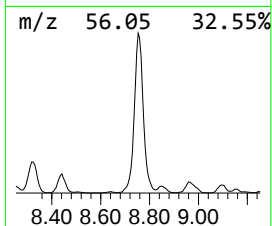
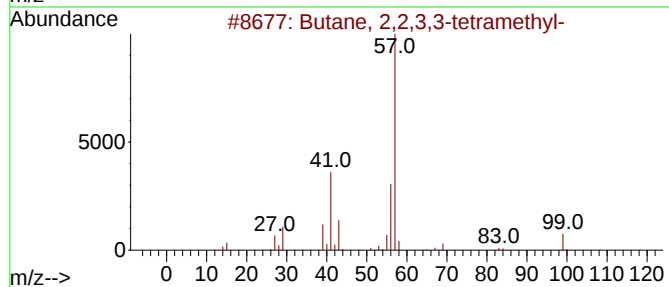
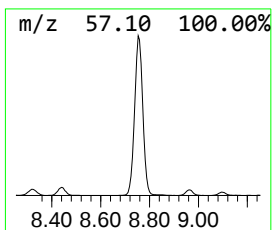
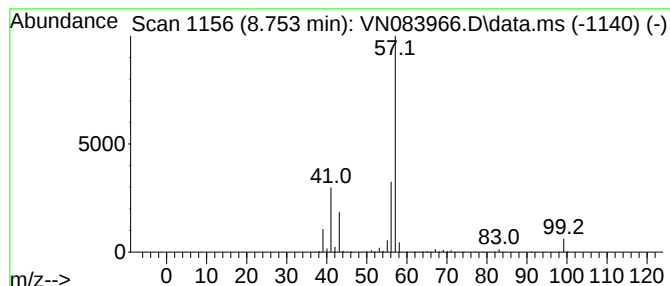
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

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

Peak Number 2 Butane, 2,2,3,3-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.753	176.49 ug/l	4219600	1,4-Difluorobenzene	9.094

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	78
2			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
3			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
4			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	64
5			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59



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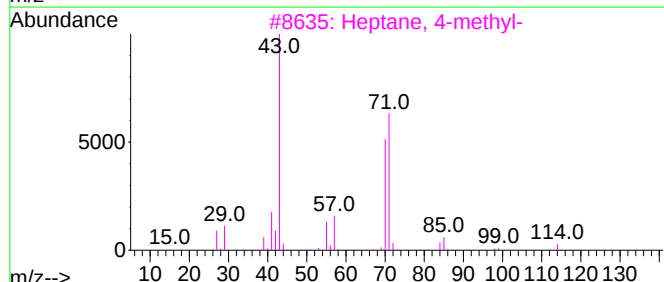
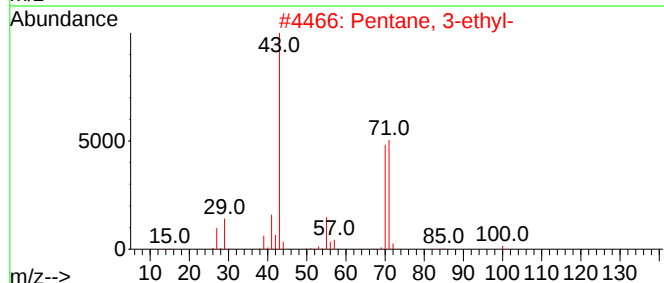
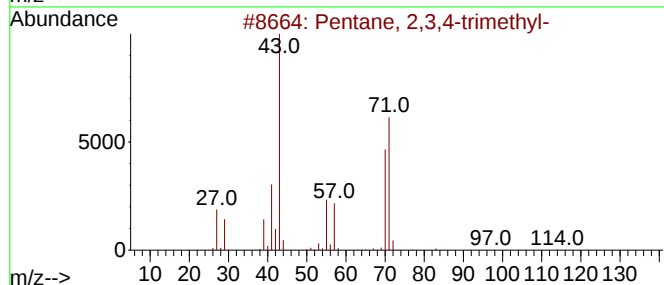
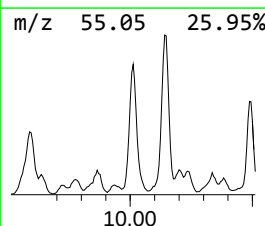
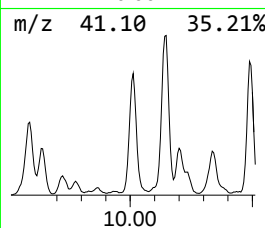
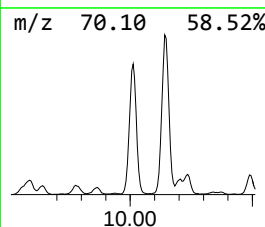
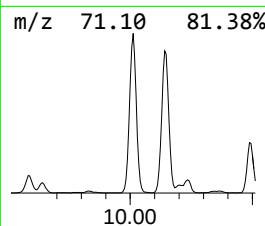
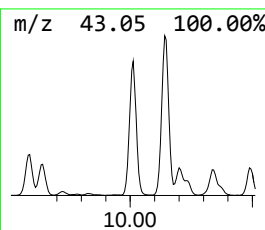
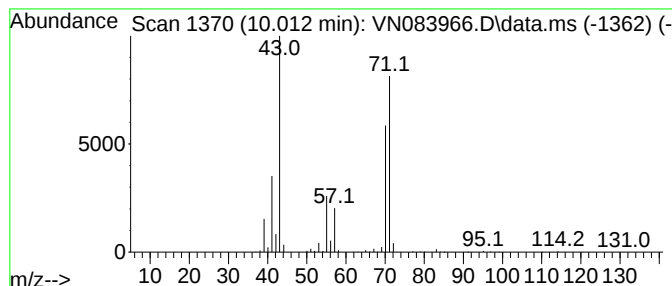
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,3,4-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.012	101.64 ug/l	2430100	1,4-Difluorobenzene	9.094

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2			Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	83
4			Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	78
5			Hexane, 3-methyl-	100	C7H16	000589-34-4	78



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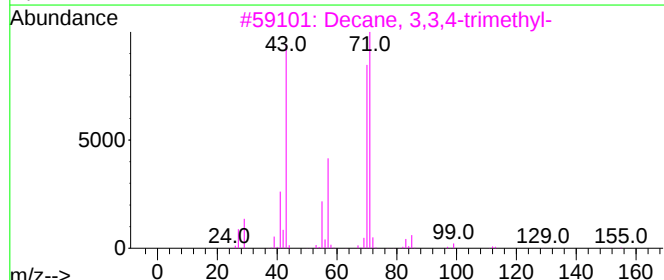
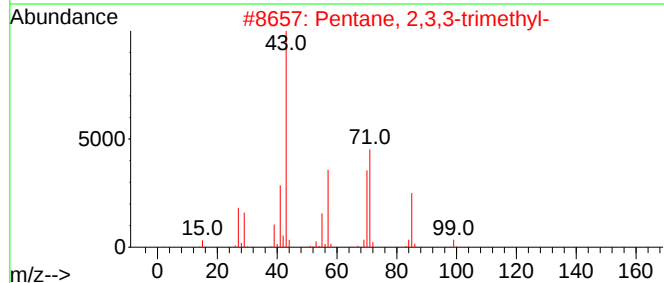
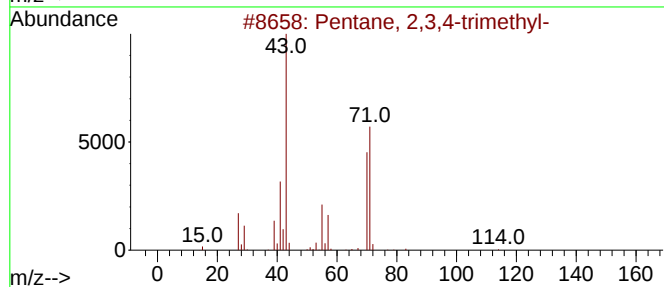
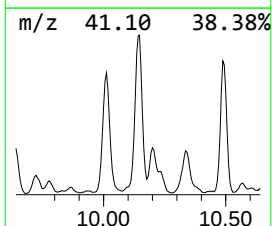
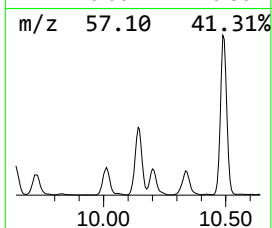
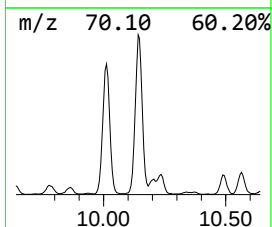
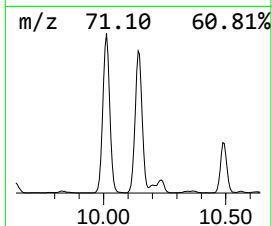
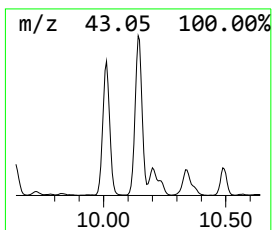
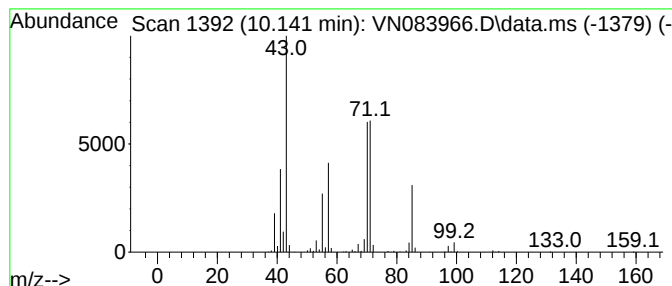
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Quant Title : SW846 8260

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

Peak Number 4 Pentane, 2,3,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.141	151.22 ug/l	3615510	1,4-Difluorobenzene	9.094

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	81
2			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	64
3			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64
4			Heptane, 4-methyl-	114	C8H18	000589-53-7	64
5			Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083966.D
Acq On : 18 Sep 2024 06:45
Operator : JC\MD
Sample : P3964-05
Misc : 10.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-2

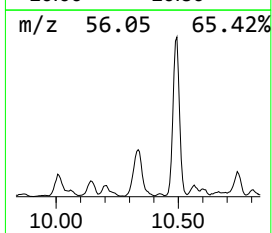
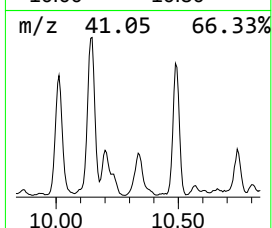
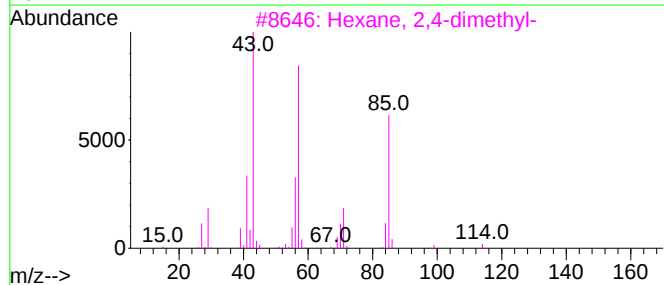
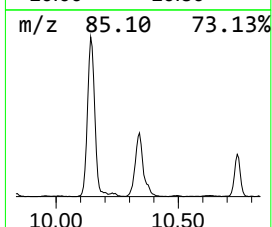
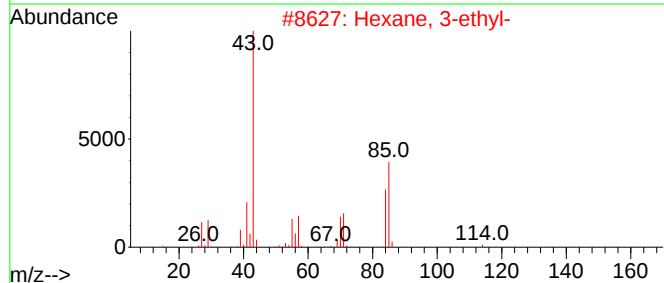
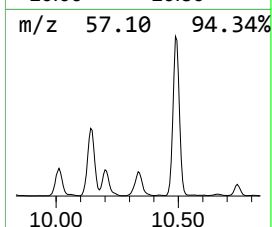
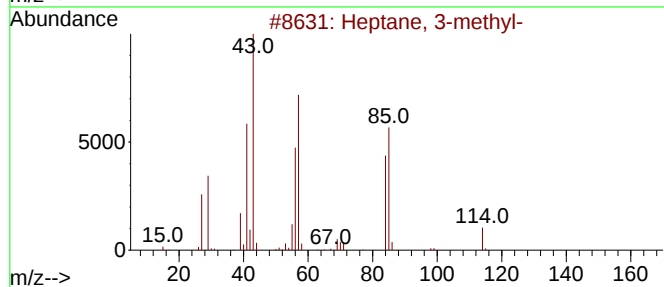
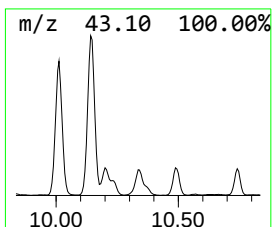
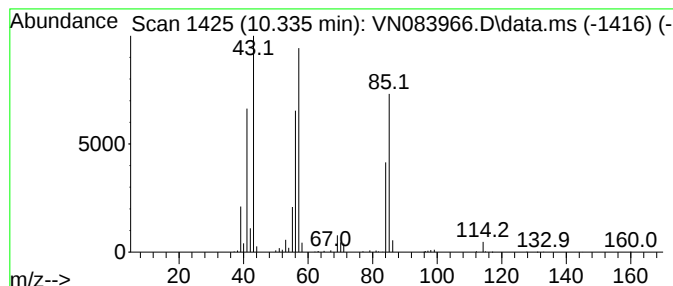
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Quant Title : SW846 8260

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

Peak Number 5 Heptane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.335	36.13 ug/l	863777	1,4-Difluorobenzene	9.094

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-methyl-	114	C8H18	000589-81-1	90
2			Hexane, 3-ethyl-	114	C8H18	000619-99-8	59
3			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	59
4			Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	59
5			Sulfurous acid, butyl isohexyl e...	222	C10H22O3S	1000309-17-2	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083966.D
Acq On : 18 Sep 2024 06:45
Operator : JC\MD
Sample : P3964-05
Misc : 10.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-2

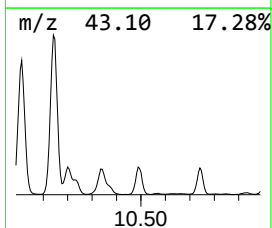
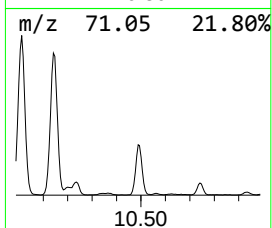
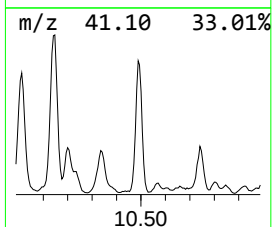
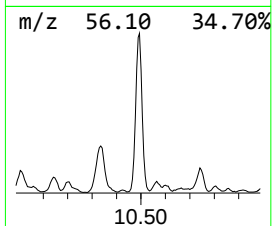
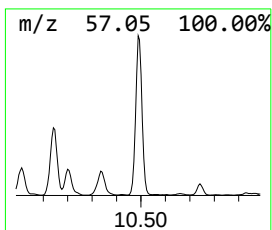
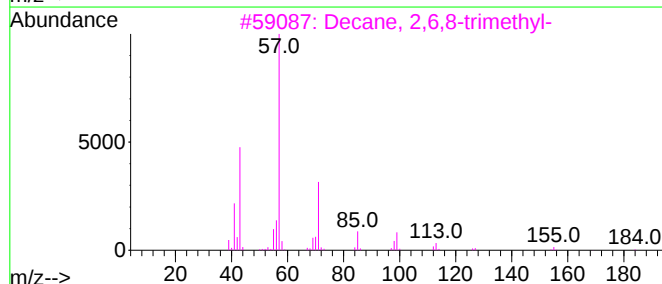
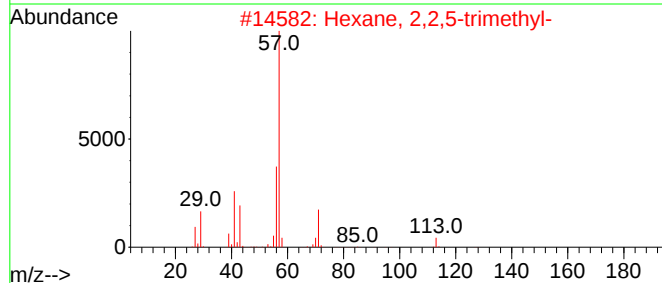
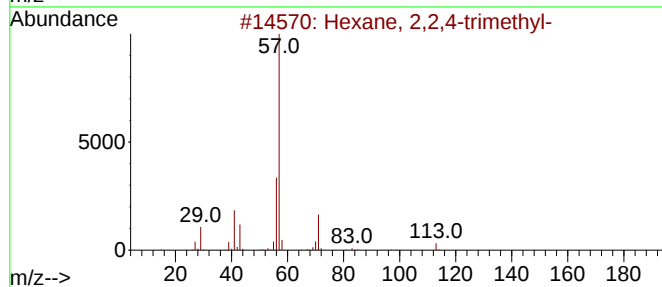
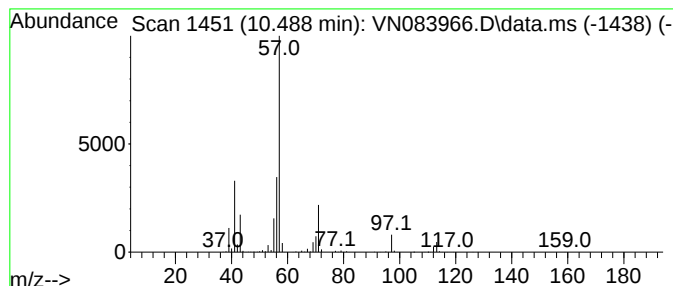
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Quant Title : SW846 8260

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

Peak Number 6 Hexane, 2,2,4-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.488	80.33 ug/l	2057250	Chlorobenzene-d5	11.865

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	72
2			Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	72
3			Decane, 2,6,8-trimethyl-	184	C13H28	062108-26-3	64
4			Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	59
5			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	56



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083966.D
Acq On : 18 Sep 2024 06:45
Operator : JC\MD
Sample : P3964-05
Misc : 10.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-2

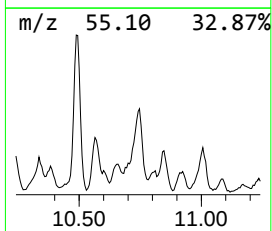
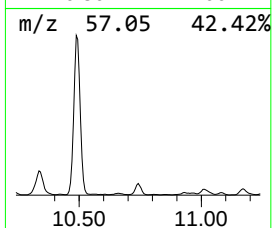
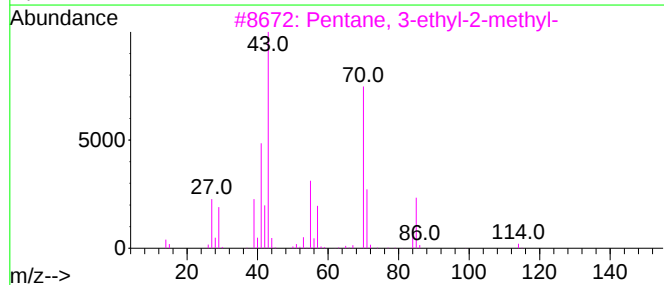
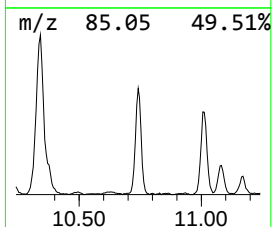
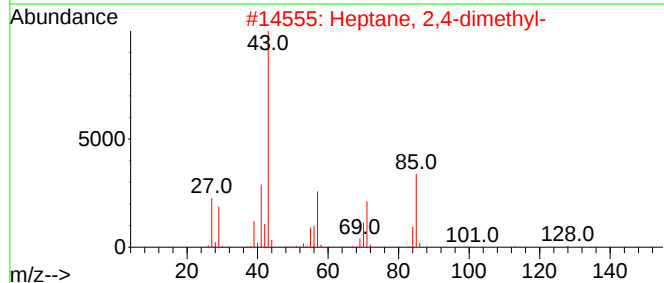
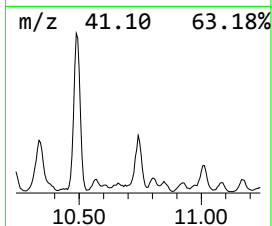
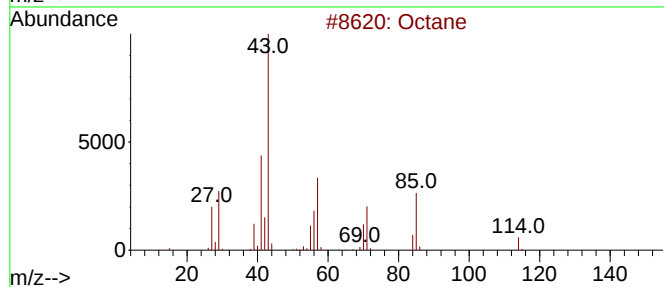
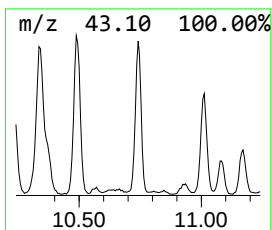
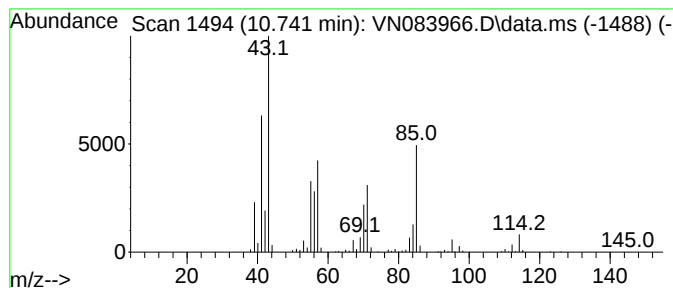
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Quant Title : SW846 8260

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

Peak Number 7 Octane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.741	25.28 ug/l	647504	Chlorobenzene-d5	11.865

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane	114	C8H18	000111-65-9	90
2			Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	64
3			Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	43
4			1-Pentanol, 2-methyl-	102	C6H14O	000105-30-6	43
5			Undecane, 2,4-dimethyl-	184	C13H28	017312-80-0	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083966.D
Acq On : 18 Sep 2024 06:45
Operator : JC\MD
Sample : P3964-05
Misc : 10.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 42 Sample Multiplier: 1

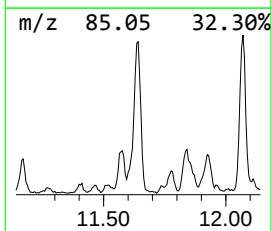
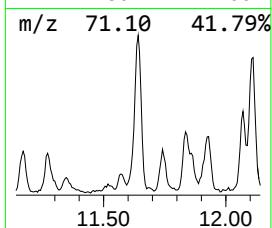
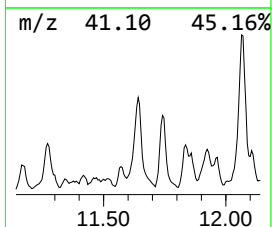
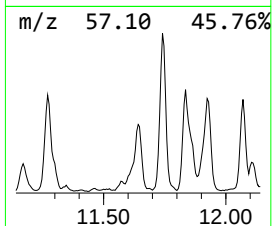
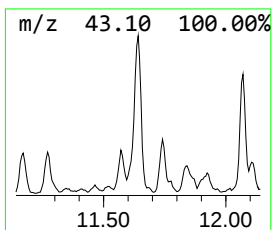
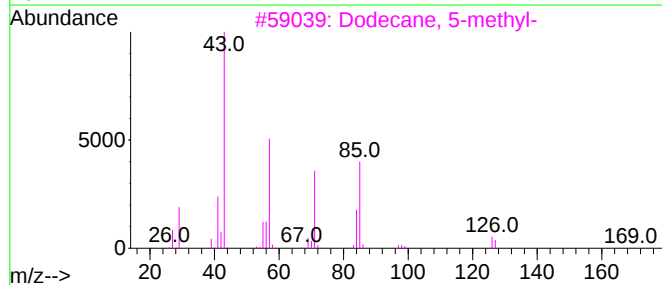
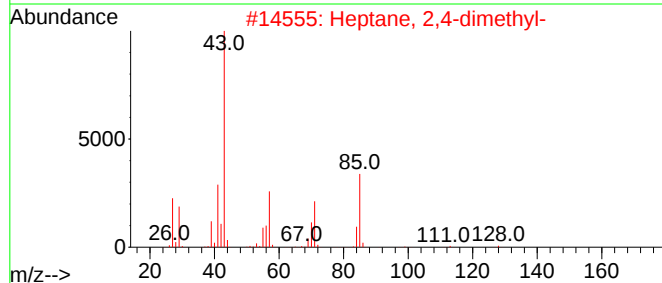
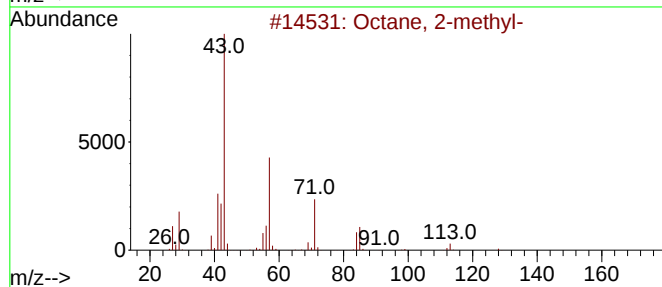
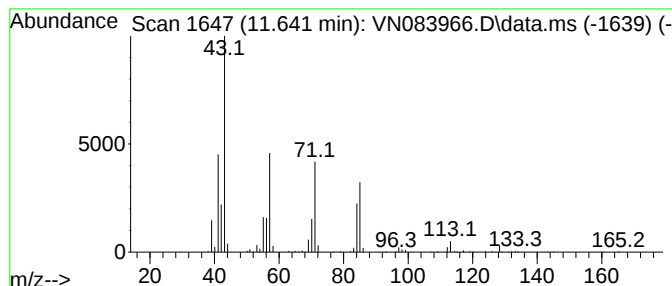
Instrument :
MSVOA_N
ClientSampleId :
SINK-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

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

Peak Number 8 Octane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.641	28.41 ug/l	727537	Chlorobenzene-d5		11.865	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octane, 2-methyl-		128	C9H20	003221-61-2	87
2	Heptane, 2,4-dimethyl-		128	C9H20	002213-23-2	64
3	Dodecane, 5-methyl-		184	C13H28	017453-93-9	59
4	Undecane, 5-methyl-		170	C12H26	001632-70-8	59
5	Undecane, 4,7-dimethyl-		184	C13H28	017301-32-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083966.D
Acq On : 18 Sep 2024 06:45
Operator : JC\MD
Sample : P3964-05
Misc : 10.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-2

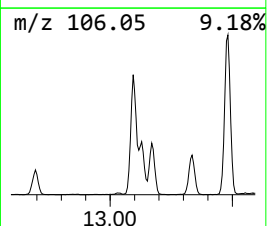
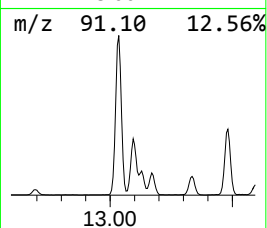
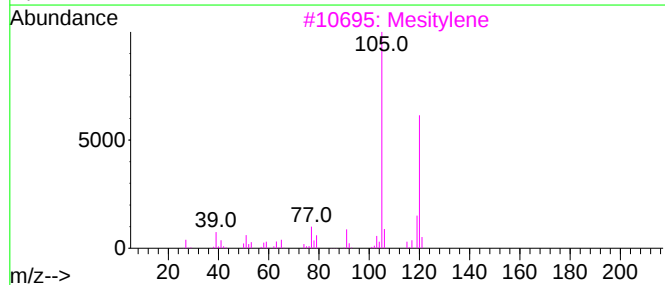
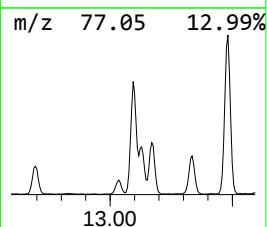
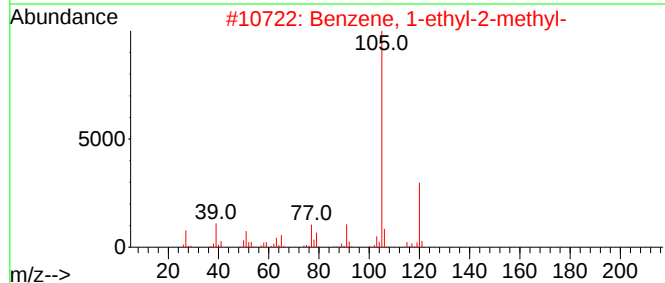
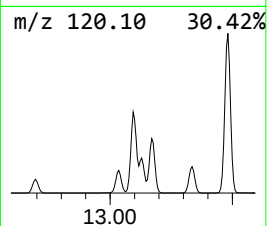
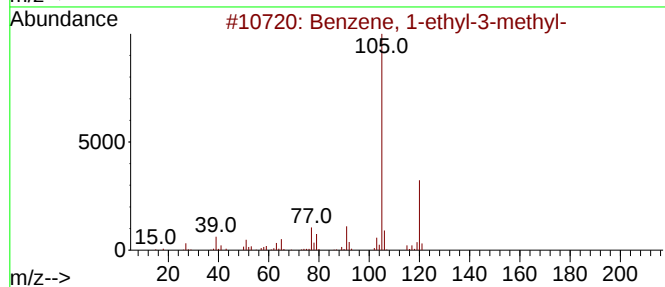
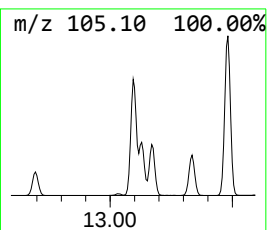
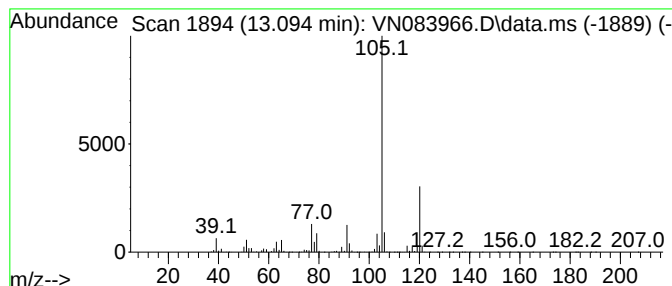
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.094	73.55 ug/l	5851090	1,4-Dichlorobenzene-d4	13.788

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083966.D
Acq On : 18 Sep 2024 06:45
Operator : JC\MD
Sample : P3964-05
Misc : 10.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-2

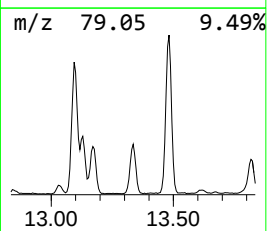
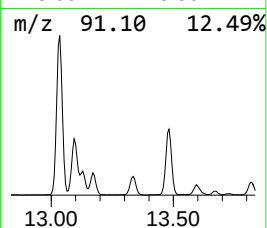
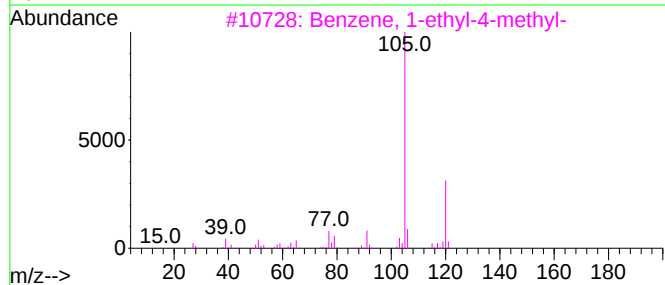
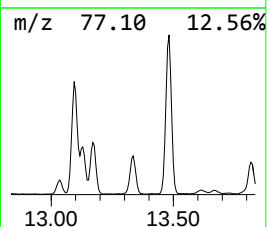
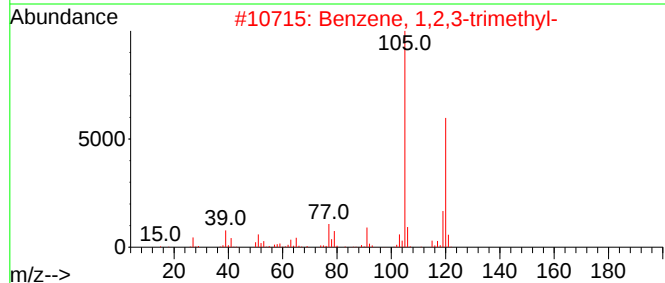
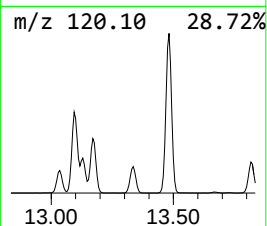
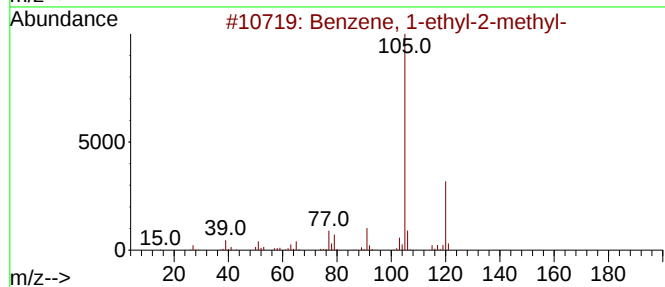
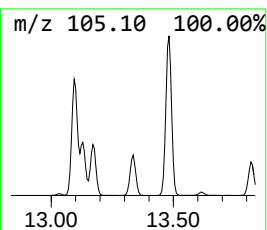
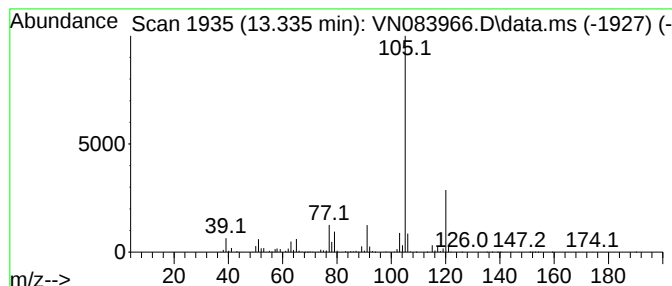
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.335	25.05 ug/l	1993170	1,4-Dichlorobenzene-d4	13.788

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083966.D
Acq On : 18 Sep 2024 06:45
Operator : JC\MD
Sample : P3964-05
Misc : 10.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.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
Hexane, 3-methyl-	8.441	27.0	ug/l	711619	1	8.218	1318700	50.0
Butane, 2,2,3,3...	8.753	176.5	ug/l	4219600	2	9.094	1195440	50.0
Pentane, 2,3,4-...	10.012	101.6	ug/l	2430100	2	9.094	1195440	50.0
Pentane, 2,3,3-...	10.141	151.2	ug/l	3615510	2	9.094	1195440	50.0
Heptane, 3-methyl-	10.335	36.1	ug/l	863777	2	9.094	1195440	50.0
Hexane, 2,2,4-t...	10.488	80.3	ug/l	2057250	3	11.865	1280510	50.0
Octane	10.741	25.3	ug/l	647504	3	11.865	1280510	50.0
Octane, 2-methyl-	11.641	28.4	ug/l	727537	3	11.865	1280510	50.0
Benzene, 1-ethy...	13.094	73.5	ug/l	5851090	4	13.788	3977640	50.0
Benzene, 1-ethy...	13.335	25.1	ug/l	1993170	4	13.788	3977640	50.0