

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN062520\
 Data File : VN062118.D
 Acq On : 25 Jun 2020 18:05
 Operator : JC/MD
 Sample : L3071-04
 Misc : 10.62g/5.0mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SB-B

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.683	267	283	304	rVB	767182	1669291	4.77%	0.360%
2	3.043	379	395	423	rVB3	310567	862020	2.47%	0.186%
3	3.487	518	533	554	rVB2	211249	513219	1.47%	0.111%
4	3.722	581	606	638	rBV2	288648	906153	2.59%	0.195%
5	4.461	815	836	840	rBV2	427404	1138521	3.26%	0.245%
6	4.979	966	997	1029	rBV	515563	1871846	5.35%	0.404%
7	5.307	1077	1099	1132	rVB4	119399	420286	1.20%	0.091%
8	5.500	1132	1159	1190	rBV	453967	1432606	4.10%	0.309%
9	5.860	1259	1271	1294	rVB	346794	1065940	3.05%	0.230%
10	6.005	1295	1316	1332	rBV2	184547	593508	1.70%	0.128%
11	6.307	1389	1410	1434	rBV4	322687	1061173	3.04%	0.229%
12	6.461	1437	1458	1475	rBV2	968295	3055627	8.74%	0.659%
13	6.567	1476	1491	1508	rVB2	651329	1913889	5.47%	0.413%
14	7.275	1691	1711	1718	rBV3	141248	375792	1.07%	0.081%
15	7.339	1719	1731	1749	rVB	446980	1198456	3.43%	0.258%
16	7.587	1773	1808	1828	rBV3	3236001	9612342	27.49%	2.072%
17	7.699	1829	1843	1864	rVB2	1621058	4411634	12.62%	0.951%
18	7.831	1864	1884	1916	rVB	3628950	8769306	25.08%	1.890%
19	7.989	1917	1933	1955	rBV3	247401	743279	2.13%	0.160%
20	8.165	1956	1988	2010	rBV	11812739	34962328	100.00%	7.537%
21	8.262	2011	2018	2039	rVB3	329238	878641	2.51%	0.189%
22	8.397	2039	2060	2086	rVB	3505495	7698841	22.02%	1.660%
23	8.542	2087	2105	2130	rBV3	1523749	4415404	12.63%	0.952%
24	8.709	2147	2157	2167	rVB	238419	451883	1.29%	0.097%
25	8.783	2167	2180	2188	rBV3	509907	1154700	3.30%	0.249%
26	9.050	2234	2263	2273	rBV3	5610545	13149796	37.61%	2.835%
27	9.108	2273	2281	2294	rVV	4022615	7745955	22.16%	1.670%
28	9.185	2294	2305	2314	rVV	1174694	2370757	6.78%	0.511%
29	9.236	2315	2321	2333	rVV3	556979	1140241	3.26%	0.246%
30	9.329	2334	2350	2365	rVV4	437749	1501745	4.30%	0.324%
31	9.487	2385	2399	2420	rBV	10846136	22729616	65.01%	4.900%
32	9.622	2421	2441	2455	rBV3	13961105	32837284	93.92%	7.079%
33	9.696	2455	2464	2471	rBV	3989556	7229431	20.68%	1.558%
34	9.837	2488	2508	2530	rBV	4667624	11897316	34.03%	2.565%

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35	9.934	2530	2538	2544	rBV3	311410	518970	1.48%	0.112%
36	9.992	2544	2556	2567	rBV	7955965	14532309	41.57%	3.133%
37	10.056	2567	2576	2588	rVV2	1504027	2942487	8.42%	0.634%
38	10.124	2591	2597	2605	rVB	795917	1228681	3.51%	0.265%
39	10.255	2622	2638	2649	rVB	4743695	9110661	26.06%	1.964%
40	10.317	2650	2657	2663	rBV2	468784	744682	2.13%	0.161%
41	10.355	2664	2669	2677	rVB2	546501	796807	2.28%	0.172%
42	10.442	2690	2696	2702	rVB2	300162	382285	1.09%	0.082%
43	10.493	2702	2712	2714	rBV4	997636	1499786	4.29%	0.323%
44	10.599	2735	2745	2756	rVB5	851172	1577783	4.51%	0.340%
45	10.689	2756	2773	2783	rBV2	903238	1825225	5.22%	0.393%
46	10.792	2796	2805	2818	rVB	1374828	2442487	6.99%	0.527%
47	10.976	2856	2862	2871	rVB3	387901	592273	1.69%	0.128%
48	11.030	2872	2879	2890	rVB5	447952	819025	2.34%	0.177%
49	11.098	2891	2900	2906	rBV2	725603	1182146	3.38%	0.255%
50	11.172	2907	2923	2943	rVB3	3879693	9572274	27.38%	2.064%
51	11.271	2944	2954	2972	rBV	3003653	5544101	15.86%	1.195%
52	11.378	2974	2987	2997	rBV3	2077488	4801245	13.73%	1.035%
53	11.484	3007	3020	3036	rVB	7007198	13799369	39.47%	2.975%
54	11.603	3038	3057	3076	rBV3	9882263	22743812	65.05%	4.903%
55	11.834	3120	3129	3137	rBV3	703471	1157054	3.31%	0.249%
56	12.091	3202	3209	3216	rBV4	437458	684328	1.96%	0.148%
57	12.136	3217	3223	3233	rVB6	403246	646454	1.85%	0.139%
58	12.223	3241	3250	3260	rBV	4071022	6505882	18.61%	1.402%
59	12.358	3282	3292	3305	rVB3	3576918	7667191	21.93%	1.653%
60	12.419	3305	3311	3319	rVB2	931987	1173404	3.36%	0.253%
61	12.522	3335	3343	3347	rBV	823618	1243895	3.56%	0.268%
62	12.564	3348	3356	3367	rVV	6158226	10361478	29.64%	2.234%
63	12.632	3368	3377	3380	rVV	2374924	3416314	9.77%	0.736%
64	12.664	3381	3387	3393	rVV	4850734	7715844	22.07%	1.663%
65	12.705	3393	3400	3421	rVB2	7257815	12874378	36.82%	2.775%
66	12.866	3440	3450	3459	rBV	1998293	3584209	10.25%	0.773%
67	12.940	3467	3473	3474	rBV2	429034	423044	1.21%	0.091%
68	12.966	3476	3481	3487	rVV	1294820	1676741	4.80%	0.361%
69	13.014	3487	3496	3507	rVB	9101993	14450053	41.33%	3.115%
70	13.072	3507	3514	3523	rVB2	913721	1472089	4.21%	0.317%
71	13.178	3524	3547	3555	rBV3	2543163	6990470	19.99%	1.507%

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72	13.352	3579	3601	3628	rVB4	3110309	14141839	40.45%	3.049%
73	13.487	3634	3643	3646	rBV	1818616	2521600	7.21%	0.544%
74	13.519	3646	3653	3659	rVV2	4539538	8026495	22.96%	1.730%
75	13.561	3660	3666	3685	rVV4	4412834	9792466	28.01%	2.111%
76	13.648	3687	3693	3701	rVB2	918815	1238563	3.54%	0.267%
77	13.715	3707	3714	3724	rVB	1804839	2968833	8.49%	0.640%
78	13.776	3725	3733	3739	rBV	1644360	2672933	7.65%	0.576%
79	13.863	3751	3760	3771	rBV	4760907	7345190	21.01%	1.583%
80	13.969	3784	3793	3803	rVB	1994797	3205002	9.17%	0.691%
81	14.101	3828	3834	3841	rBV2	376518	533163	1.52%	0.115%
82	14.185	3851	3860	3866	rBV	2445226	3719199	10.64%	0.802%
83	14.223	3866	3872	3880	rVB	2213115	3066788	8.77%	0.661%
84	14.268	3880	3886	3893	rBV3	1280438	1751878	5.01%	0.378%
85	14.310	3894	3899	3903	rVV	551416	676162	1.93%	0.146%
86	14.342	3904	3909	3922	rVB2	1018892	1533656	4.39%	0.331%
87	14.410	3923	3930	3935	rBV2	672052	874730	2.50%	0.189%
88	14.451	3935	3943	3952	rVV2	1517365	2681880	7.67%	0.578%
89	14.500	3953	3958	3965	rVB	1102511	1432272	4.10%	0.309%
90	14.554	3965	3975	3990	rBV5	3587399	8397318	24.02%	1.810%
91	14.625	3990	3997	4007	rVB	997977	1628175	4.66%	0.351%
92	14.828	4051	4060	4080	rVV4	1402253	3550591	10.16%	0.765%
93	14.930	4083	4092	4106	rVB2	989219	2005676	5.74%	0.432%
94	15.104	4134	4146	4157	rVB	1561482	3013343	8.62%	0.650%
95	15.213	4171	4180	4188	rBV3	543274	960369	2.75%	0.207%
96	15.387	4223	4234	4248	rBV	626293	1262834	3.61%	0.272%
97	15.545	4267	4283	4289	rBV5	352689	961170	2.75%	0.207%
98	15.664	4312	4320	4329	rBV	276063	500110	1.43%	0.108%
99	16.123	4453	4463	4475	rVB	894645	1618779	4.63%	0.349%
100	16.310	4510	4521	4548	rVB	545718	1320766	3.78%	0.285%

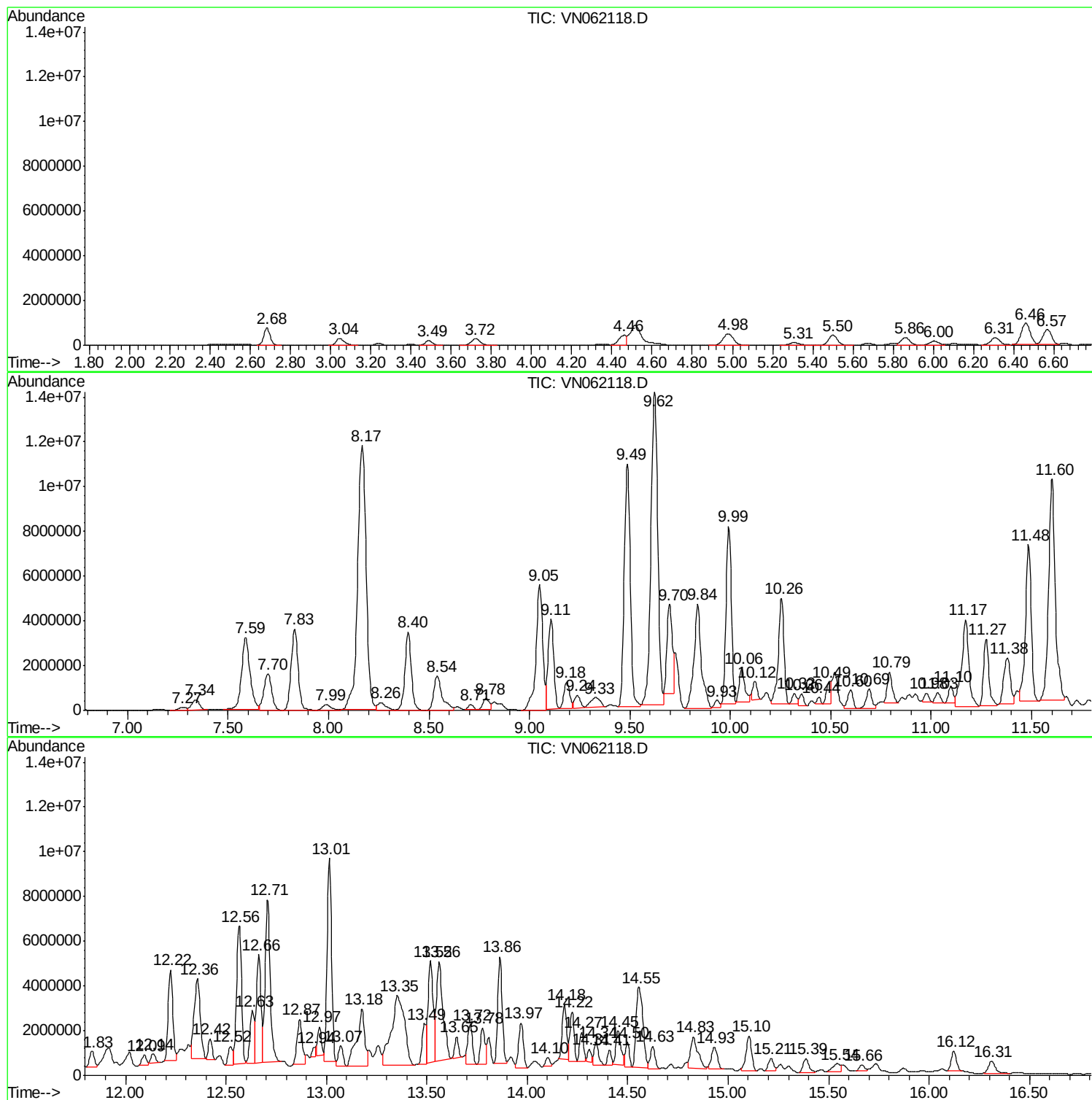
Sum of corrected areas: 463879842

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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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



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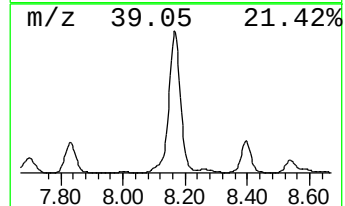
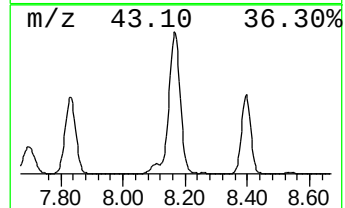
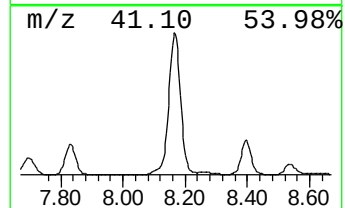
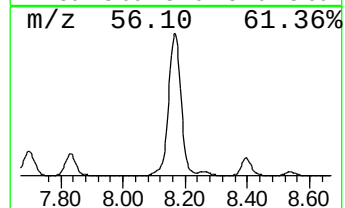
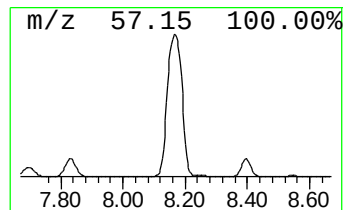
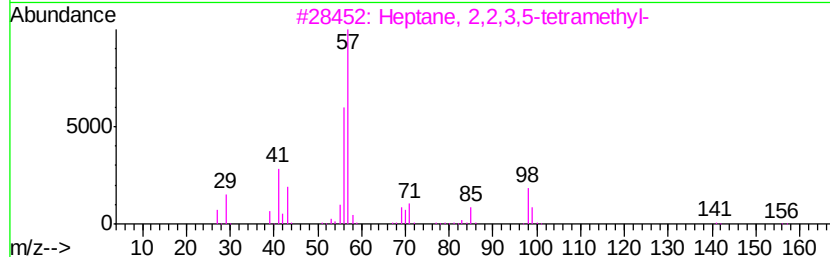
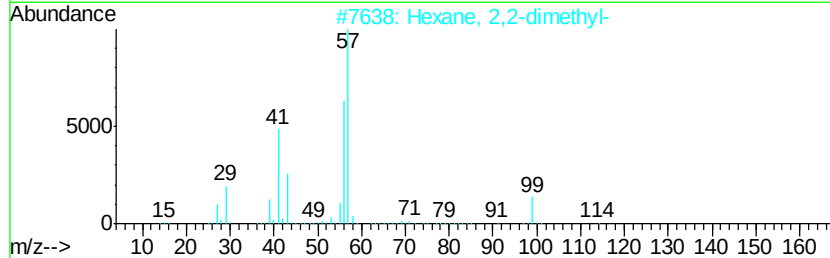
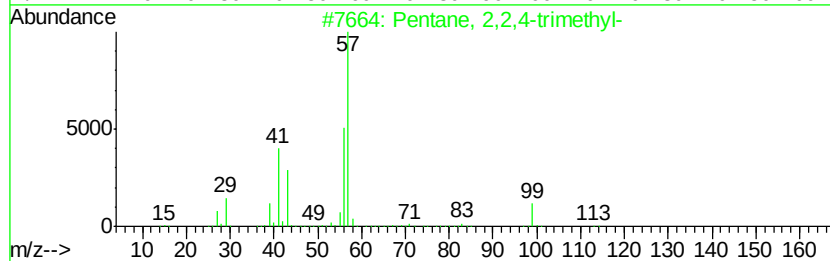
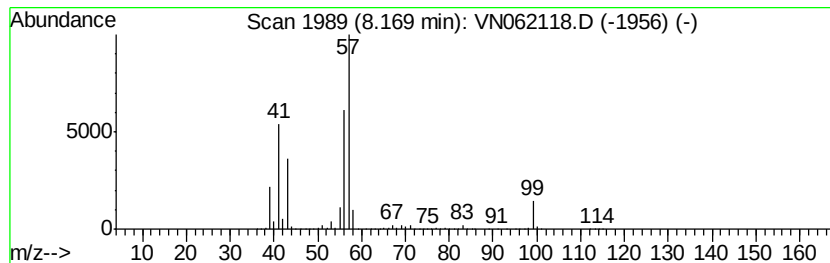
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2,2,4-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.17	395.91 ug/l	34962300	1,4-Difluorobenzene	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	83
2		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
3		Heptane, 2,2,3,5-tetramethyl-	156	C11H24	061868-42-6	64
4		Pentane, 3-methyl-	86	C6H14	000096-14-0	59
5		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	50



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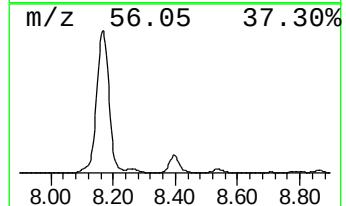
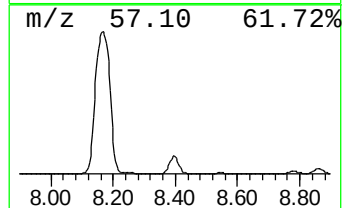
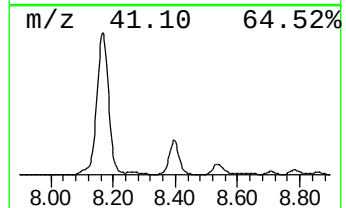
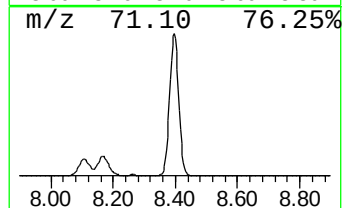
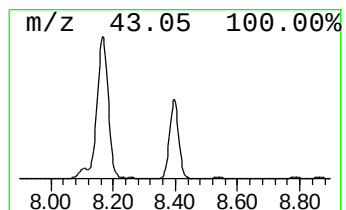
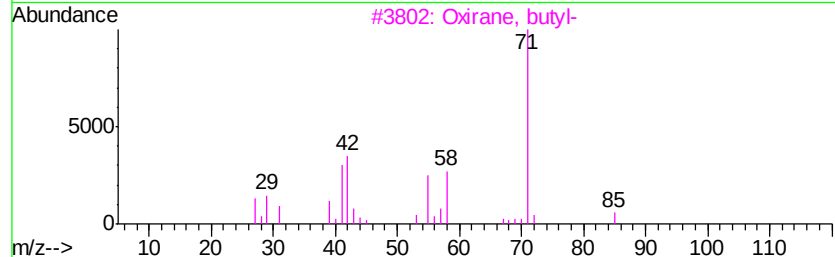
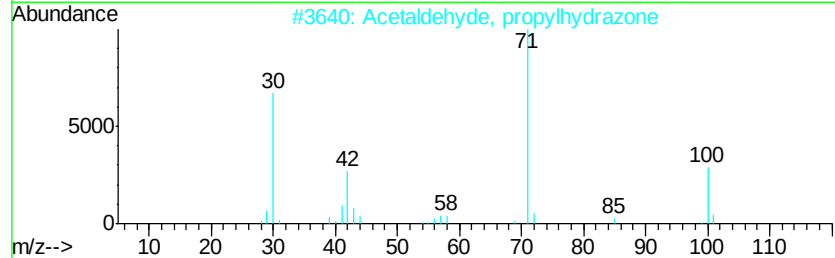
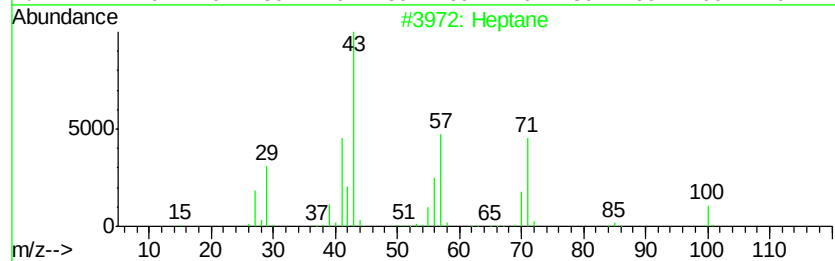
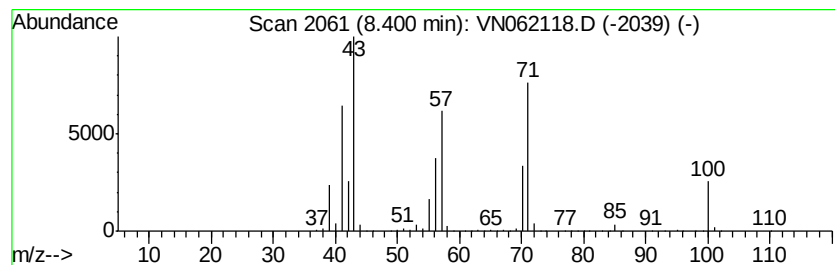
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Quant Title : SW846 8260

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

Peak Number 2 Heptane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.40	87.18 ug/l	7698840	1,4-Difluorobenzene	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	91
2		Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	58
3		Oxirane, butyl-	100	C6H12O	001436-34-6	47
4		Hexane, 3-methyl-	100	C7H16	000589-34-4	43
5		1-Hexene, 3,4,5-trimethyl-	126	C9H18	056728-10-0	42



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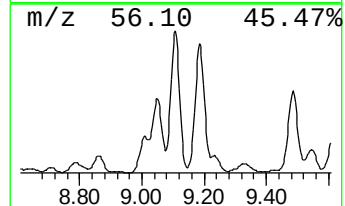
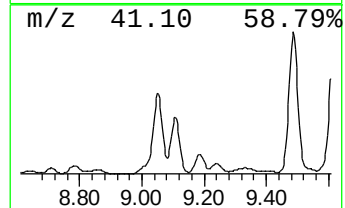
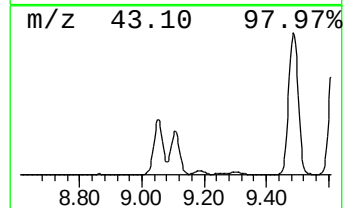
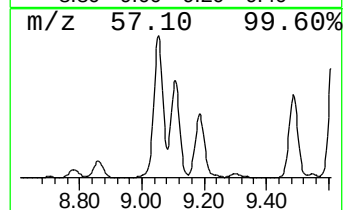
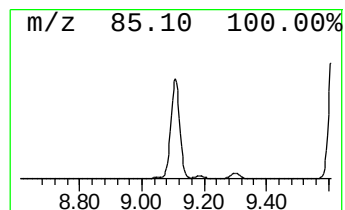
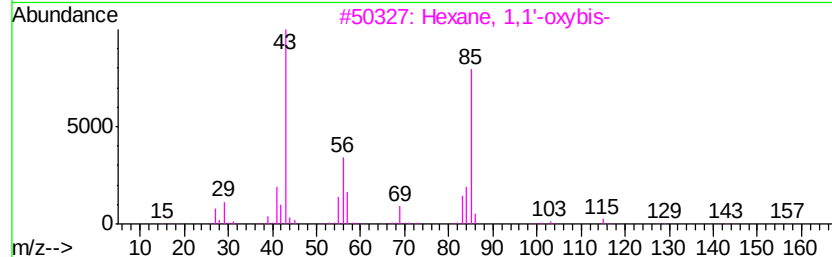
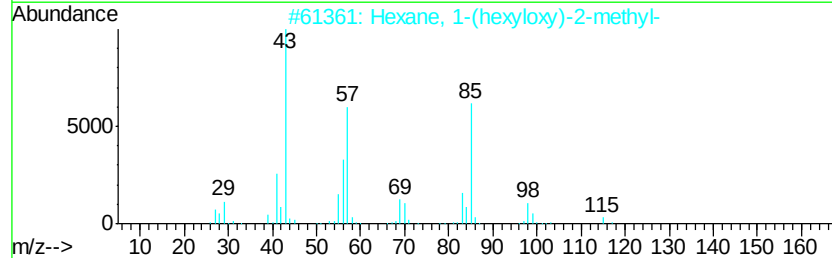
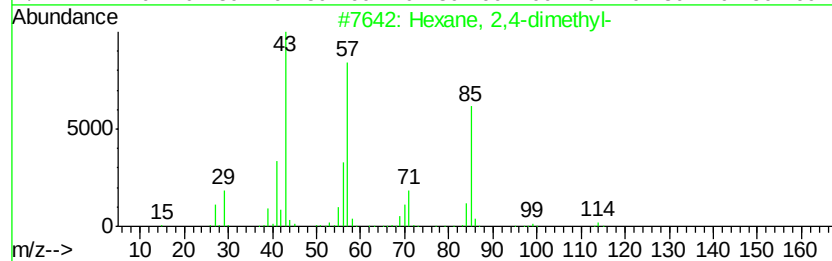
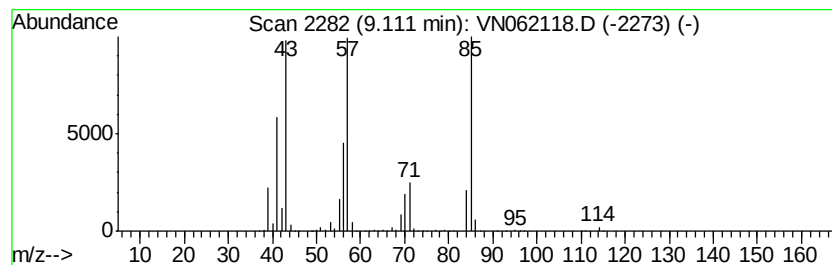
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Quant Title : SW846 8260

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

Peak Number 3 Hexane, 2,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.11	87.72 ug/l	7745960	1,4-Difluorobenzene	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	83
2		Hexane, 1-(hexyloxy)-2-methyl-	200	C13H28O	074421-17-3	72
3		Hexane, 1,1'-oxybis-	186	C12H26O	000112-58-3	64
4		Sulfurous acid, butyl hexyl ester	222	C10H22O3S	1000309-17-3	59
5		Hexane, 2-methyl-	100	C7H16	000591-76-4	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_N\DATA\VN062520\
Data File : VN062118.D
Acq On : 25 Jun 2020 18:05
Operator : JC/MD
Sample : L3071-04
Misc : 10.62µL/5.0mL/100µL/5.00mL/MSVOA_N/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SB-B

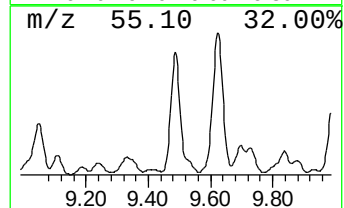
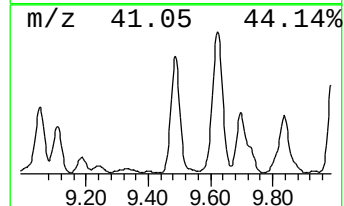
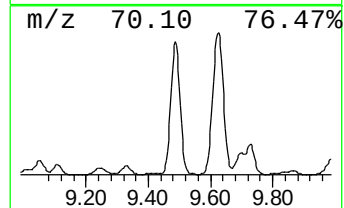
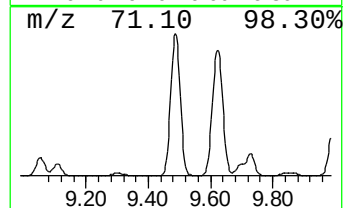
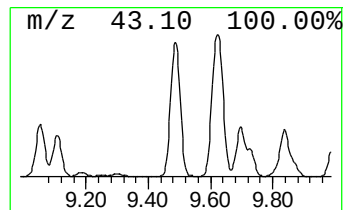
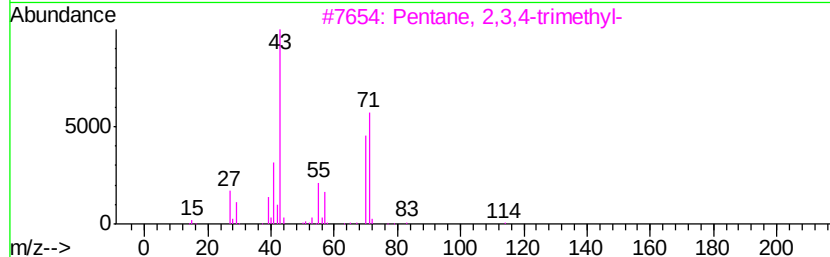
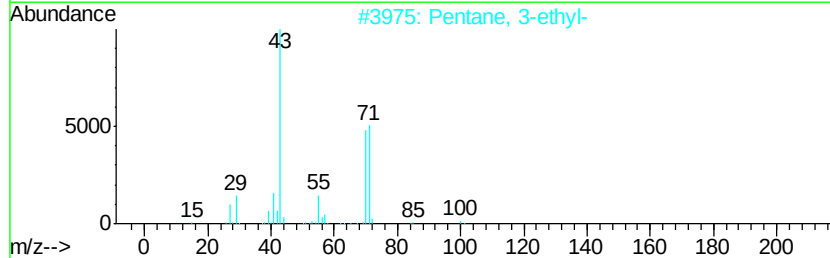
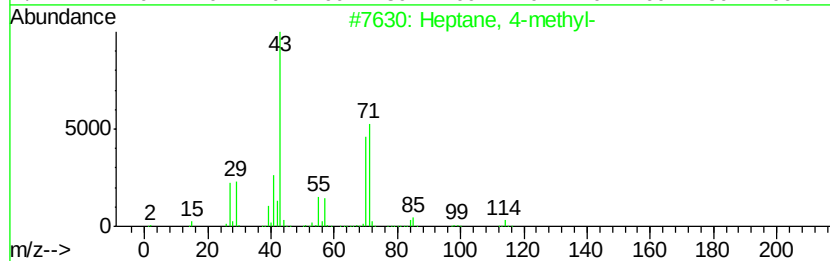
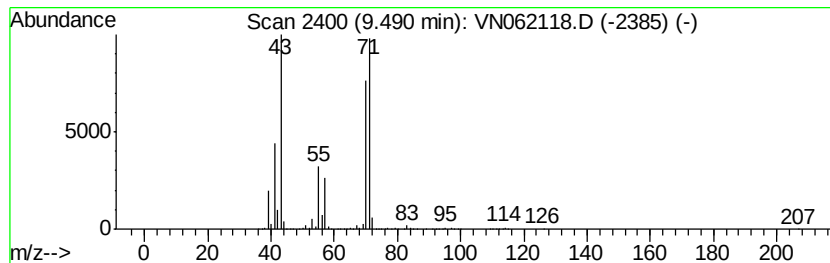
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.49	257.39 ug/l	22729600	1,4-Difluorobenzene	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane, 4-methyl-	114	C8H18	000589-53-7	91
2		Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
3		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	83
4		Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	83
5		Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_N\DATA\VN062520\
Data File : VN062118.D
Acq On : 25 Jun 2020 18:05
Operator : JC/MD
Sample : L3071-04
Misc : 10.62µ/5.0mL/100µL/5.00mL/MSVOA_N/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SB-B

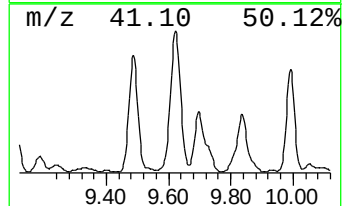
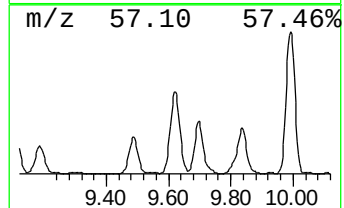
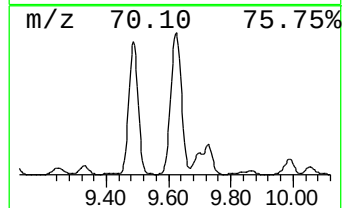
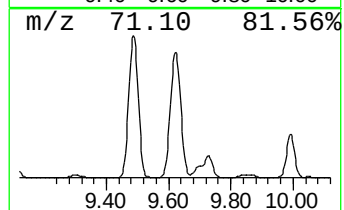
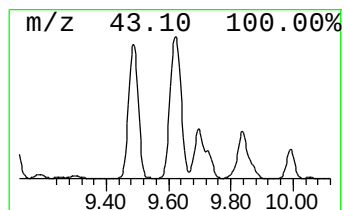
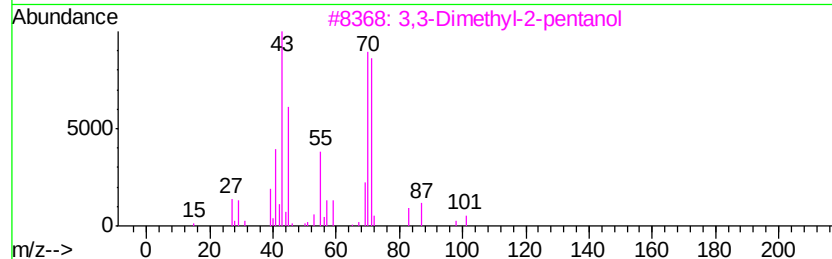
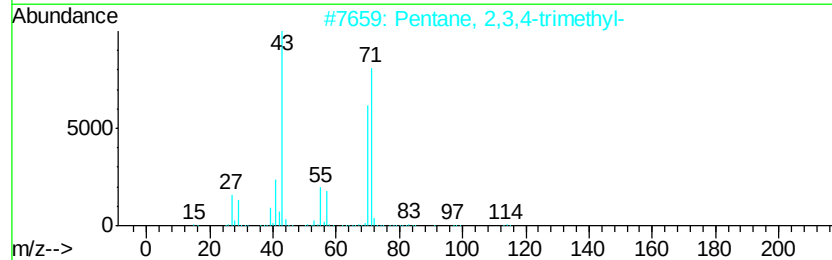
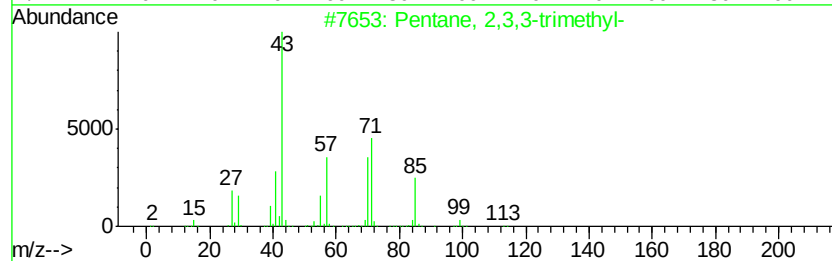
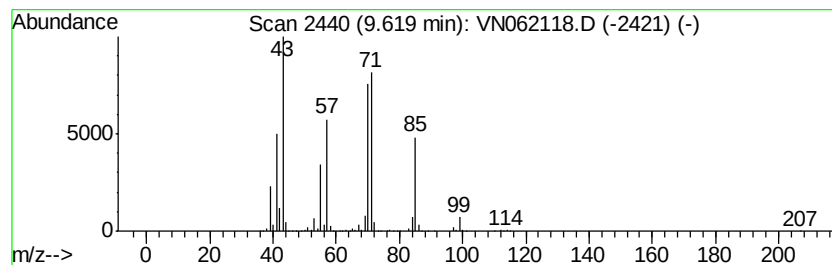
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Quant Title : SW846 8260

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

Peak Number 5 Pentane, 2,3,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.62	371.85 ug/l	32837300	1,4-Difluorobenzene	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	78
2		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	72
3		3,3-Dimethyl-2-pentanol	116	C7H16O	019781-24-9	64
4		Hexane, 3-methyl-	100	C7H16	000589-34-4	64
5		Pentane, 3-ethyl-	100	C7H16	000617-78-7	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062520\
Data File : VN062118.D
Acq On : 25 Jun 2020 18:05
Operator : JC/MD
Sample : L3071-04
Misc : 10.62uL/5.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SB-B

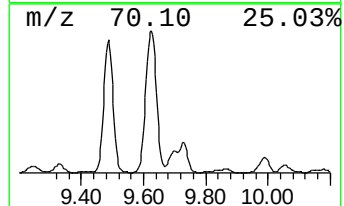
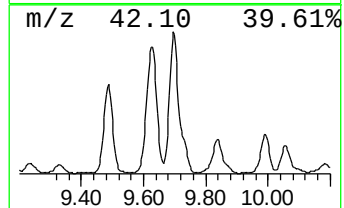
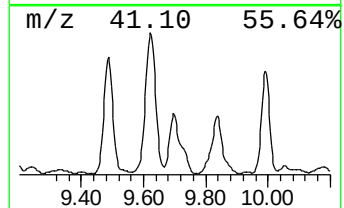
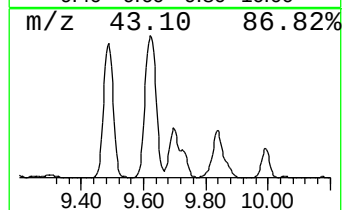
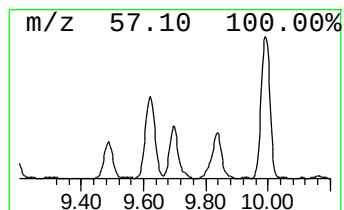
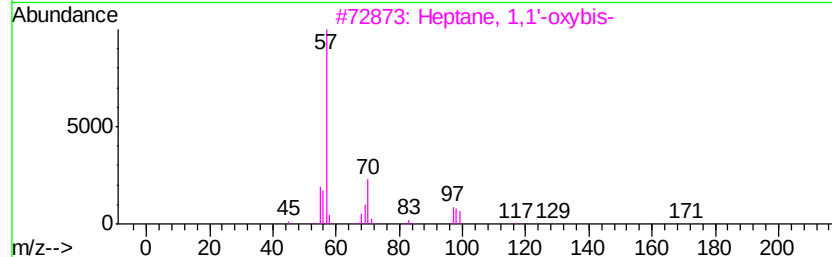
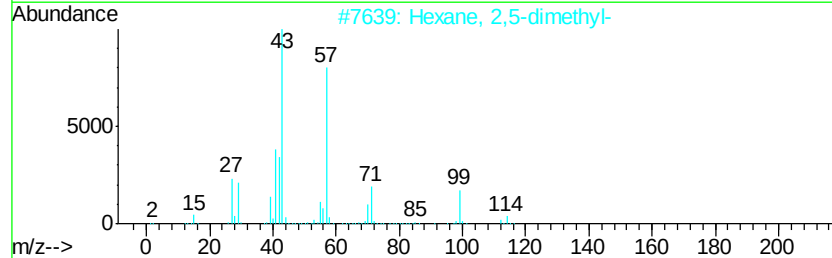
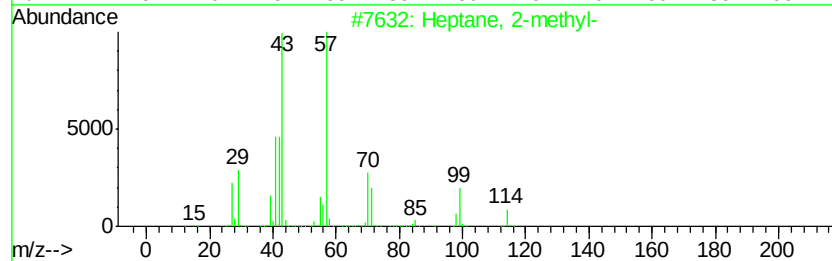
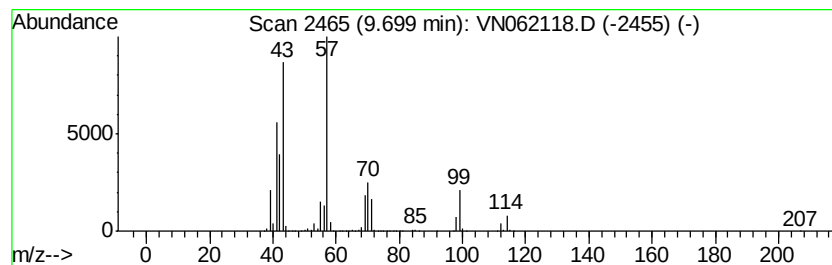
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Quant Title : SW846 8260

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

Peak Number 6 Heptane, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.70	81.87 ug/l	7229430	1,4-Difluorobenzene	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane, 2-methyl-	114	C8H18	000592-27-8	90
2		Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	64
3		Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	42
4		Heptane, 3-ethyl-	128	C9H20	015869-80-4	37
5		4-Bromoheptane	178	C7H15Br	000998-93-6	35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_N\DATA\VN062520\
Data File : VN062118.D
Acq On : 25 Jun 2020 18:05
Operator : JC/MD
Sample : L3071-04
Misc : 10.62µL/5.0mL/100µL/5.00mL/MSVOA_N/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SB-B

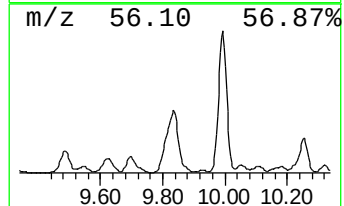
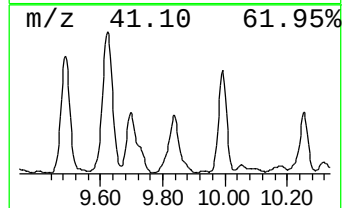
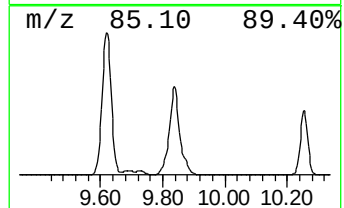
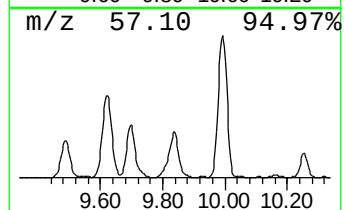
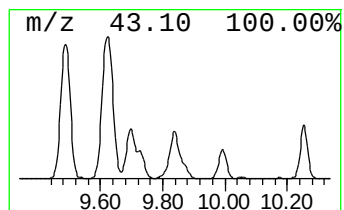
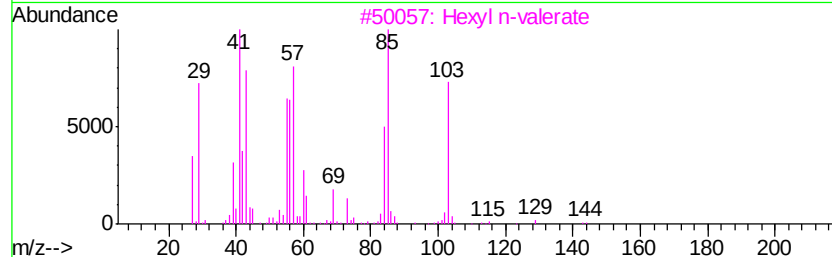
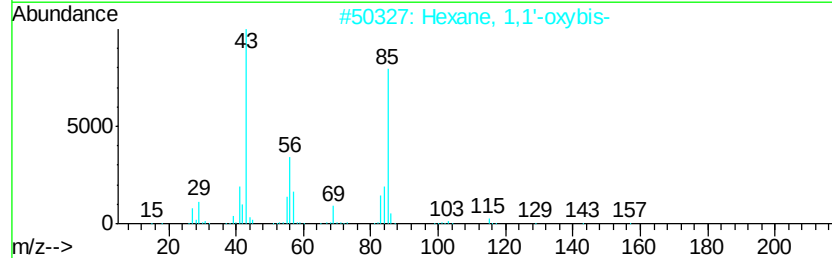
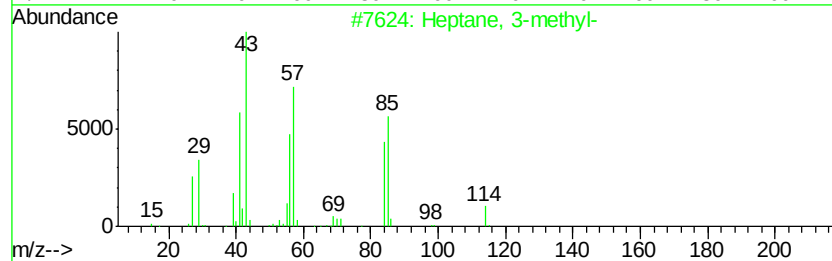
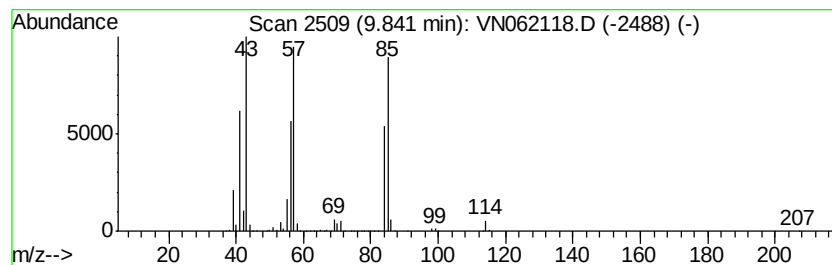
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Quant Title : SW846 8260

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

Peak Number 7 Heptane, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.84	134.73 ug/l	11897300	1,4-Difluorobenzene	8.55

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 3-methyl-	114	C8H18	000589-81-1	91
2		Hexane, 1,1'-oxybis-	186	C12H26O	000112-58-3	64
3		Hexyl n-valerate	186	C11H22O2	001117-59-5	64
4		Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	59
5		Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_N\DATA\VN062520\
Data File : VN062118.D
Acq On : 25 Jun 2020 18:05
Operator : JC/MD
Sample : L3071-04
Misc : 10.62µL/5.0mL/100µL/5.00mL/MSVOA_N/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SB-B

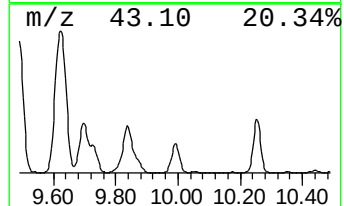
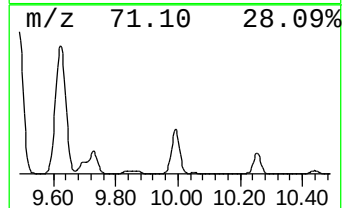
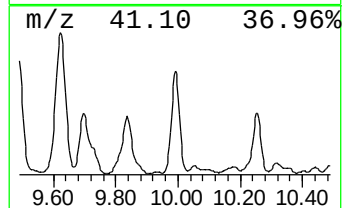
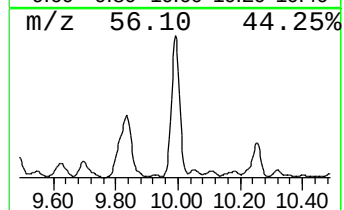
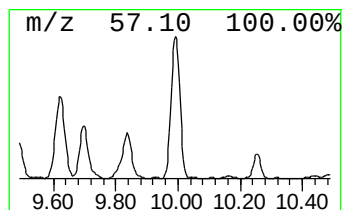
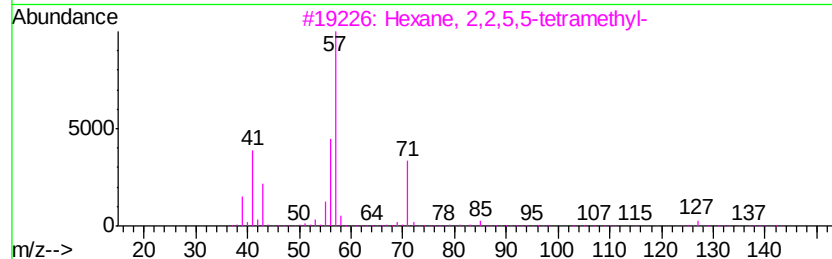
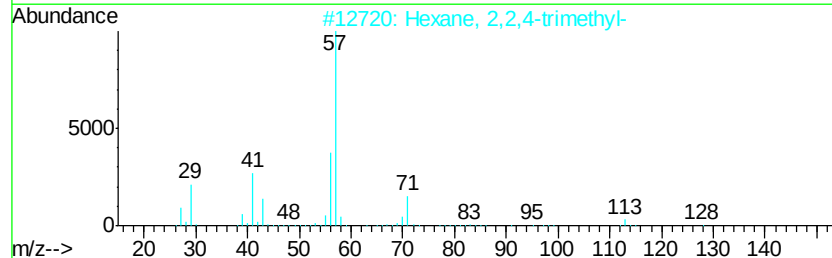
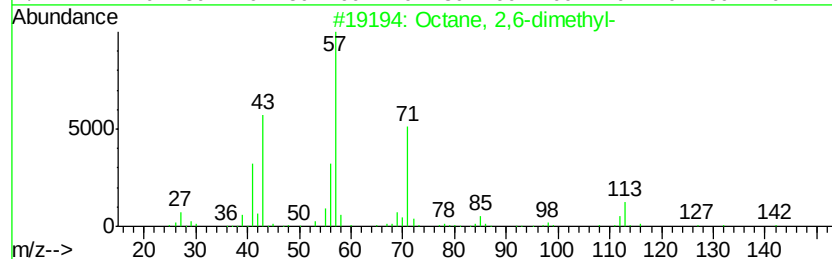
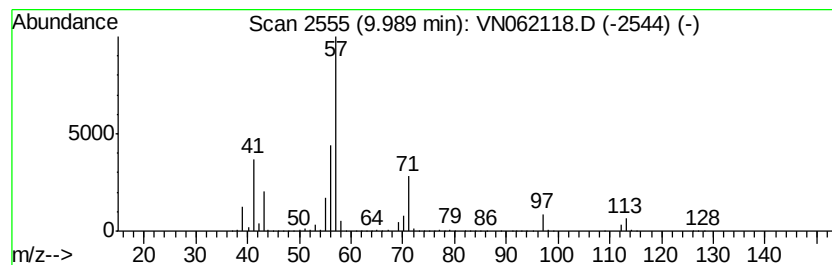
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Quant Title : SW846 8260

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

Peak Number 8 Octane, 2,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.99	151.34 ug/l	14532300	Chlorobenzene-d5	11.38

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
2		Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	53
3		Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	53
4		Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	53
5		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_N\DATA\VN062520\
Data File : VN062118.D
Acq On : 25 Jun 2020 18:05
Operator : JC/MD
Sample : L3071-04
Misc : 10.62µL/5.0mL/100µL/5.00mL/MSVOA_N/MEOH
ALS Vial : 20 Sample Multiplier: 1

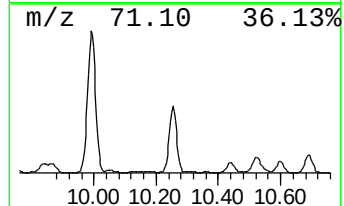
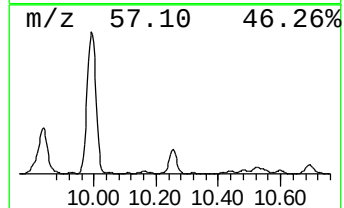
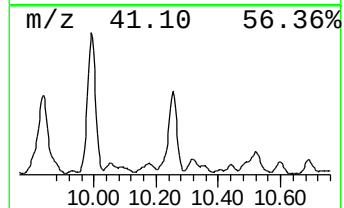
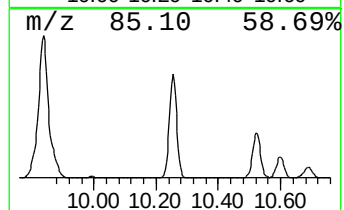
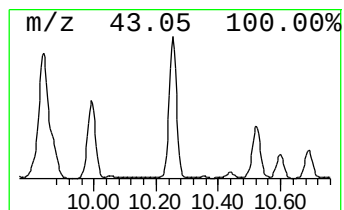
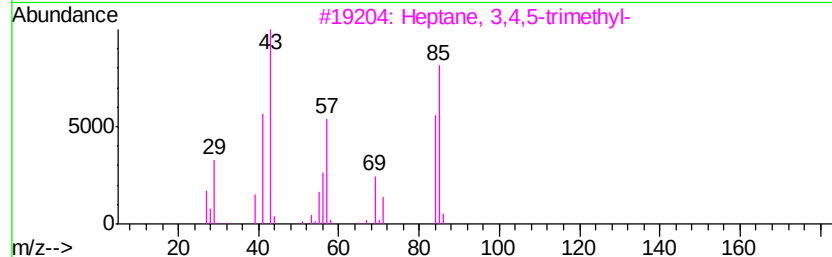
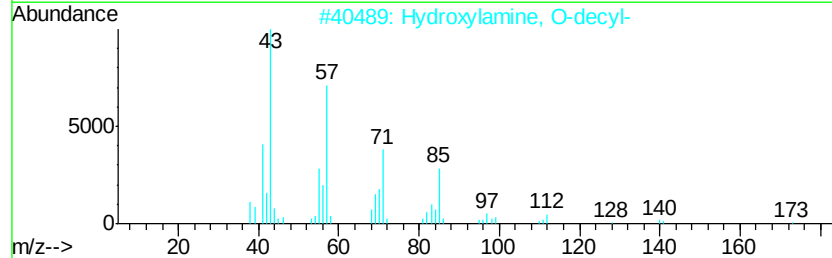
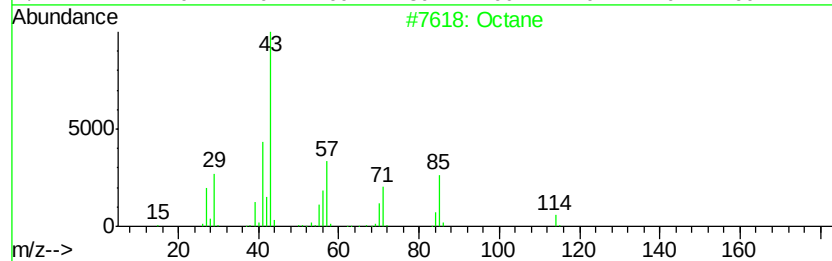
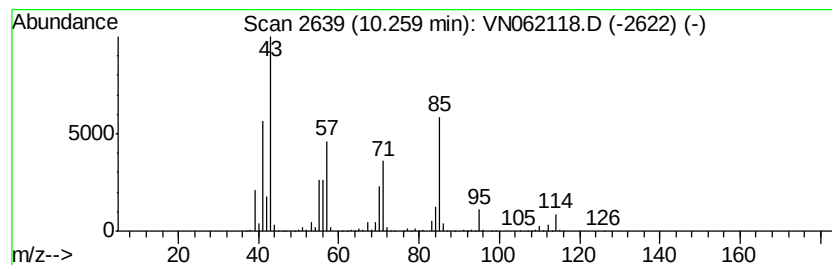
Instrument :
MSVOA_N
ClientSampleId :
SB-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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

Peak Number 9 Octane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.26	94.88 ug/l	9110660	Chlorobenzene-d5	11.38
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octane		114 C8H18	000111-65-9 91
2	Hydroxylamine, O-decyl-		173 C10H23NO	029812-79-1 59
3	Heptane, 3,4,5-trimethyl-		142 C10H22	020278-89-1 53
4	Hexane, 2-methyl-		100 C7H16	000591-76-4 53
5	Heptane, 4,4-dimethyl-		128 C9H20	001068-19-5 50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062520\
Data File : VN062118.D
Acq On : 25 Jun 2020 18:05
Operator : JC/MD
Sample : L3071-04
Misc : 10.62µ/5.0mL/100µL/5.00mL/MSVOA_N/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
SB-B

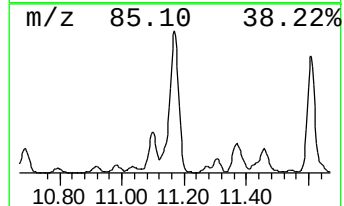
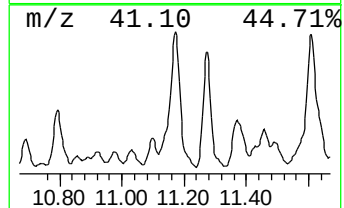
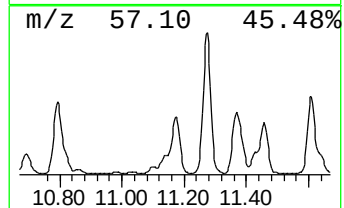
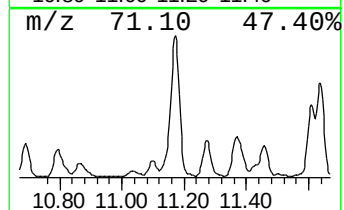
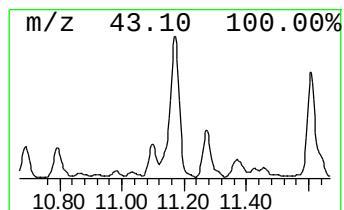
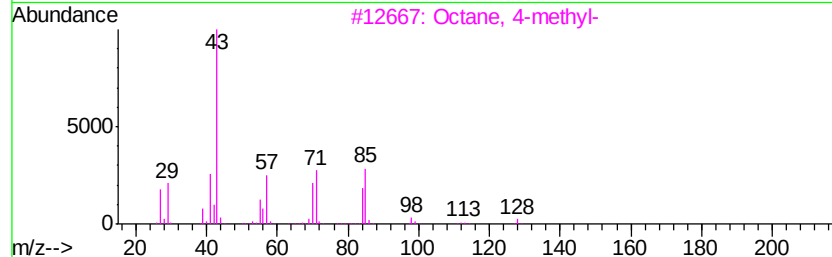
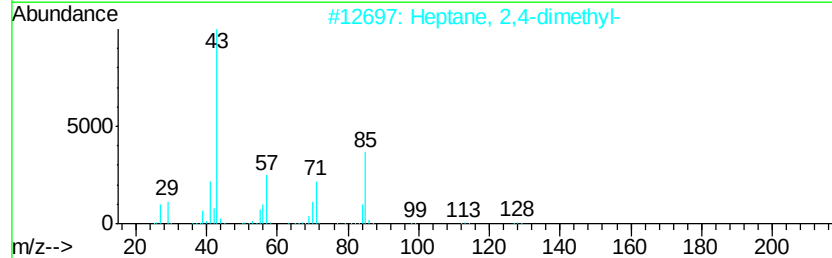
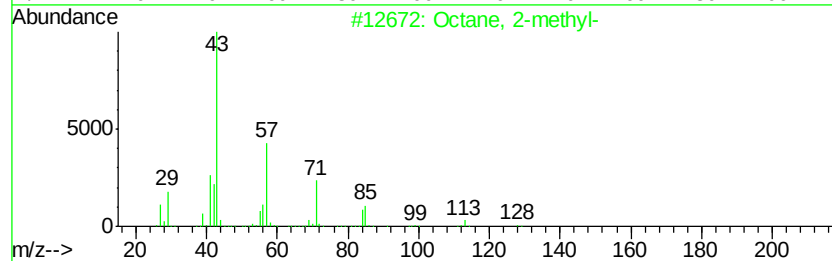
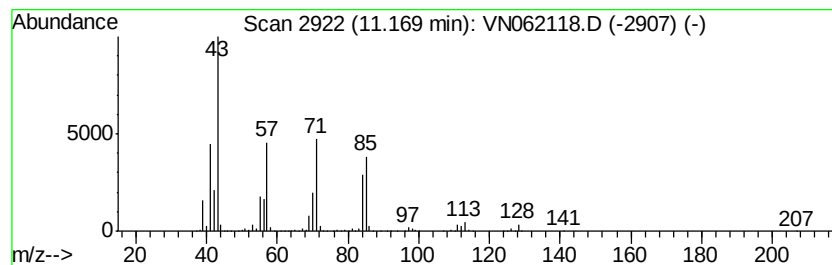
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Quant Title : SW846 8260

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

Peak Number 10 Octane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.17	99.69 ug/l	9572270	Chlorobenzene-d5	11.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2-methyl-	128	C9H20	003221-61-2	64
2		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	59
3		Octane, 4-methyl-	128	C9H20	002216-34-4	53
4		Dichloroacetic acid, 4-methylpen...	212	C8H14Cl2O2	1000282-43-7	53
5		Formic acid, 2-methylpentyl ester	130	C7H14O2	381670-34-4	50



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN062520\
Data File : VN062118.D
Acq On : 25 Jun 2020 18:05
Operator : JC/MD
Sample : L3071-04
Misc : 10.62g/5.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SB-B

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Pentane, 2,2,4-tr...	8.17	395.9	ug/l	34962300	2	8.55	4415400	50.0
Heptane	8.40	87.2	ug/l	7698840	2	8.55	4415400	50.0
Hexane, 2,4-dimet...	9.11	87.7	ug/l	7745960	2	8.55	4415400	50.0
Heptane, 4-methyl-	9.49	257.4	ug/l	22729600	2	8.55	4415400	50.0
Pentane, 2,3,3-tr...	9.62	371.9	ug/l	32837300	2	8.55	4415400	50.0
Heptane, 2-methyl-	9.70	81.9	ug/l	7229430	2	8.55	4415400	50.0
Heptane, 3-methyl-	9.84	134.7	ug/l	11897300	2	8.55	4415400	50.0
Octane, 2,6-dimet...	9.99	151.3	ug/l	14532300	3	11.38	4801250	50.0
Octane	10.26	94.9	ug/l	9110660	3	11.38	4801250	50.0
Octane, 2-methyl-	11.17	99.7	ug/l	9572270	3	11.38	4801250	50.0