LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN101019\
Data File : VN058649.D
Acq On : 9 Oct 2019 18:48
Operator : JC/SP
Sample : K5184-08
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

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\82N092719W.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	2.002	61	71	81	rBV	164943	238878	3.18%	0.188%
2	2.166	104	122	138	rBV	749917	1146753	15.26%	0.901%
3	2.789	298	316	345	rBV	3729098	7516045	100.00%	5.903%
4	3.124	400	420	449	rBV3	1187227	3120330	41.52%	2.451%
5	3.304	462	476	493	rVB	66928	145552	1.94%	0.114%
6	3.458	496	524	534	rBV4	35797	120457	1.60%	0.095%
7	3.548	534	552	584	rVB	1635452	3944449	52.48%	3.098%
8	3.780	603	624	655	rVB3	189457	595805	7.93%	0.468%
9	4.403	794	818	832	rBV3	138000	463116	6.16%	0.364%
10	4.571	833	870	933	rVB6	1161787	7378927	98.18%	5.795%
11	5.030	981	1013	1051	rBV2	940692	3408810	45.35%	2.677%
12	5.355	1083	1114	1149	rBV2	148009	516509	6.87%	0.406%
13	5.548	1149	1174	1201	rVV2	303918	963194	12.82%	0.757%
14	5.722	1207	1228	1249	rBV3	61507	194415	2.59%	0.153%
15	5.905	1262	1285	1308	rBV	656383	1999996	26.61%	1.571%
16	6.046	1308	1329	1347	rBV2	444301	1422576	18.93%	1.117%
17	6.137	1349	1357	1375	rVB2	83186	221950	2.95%	0.174%
18	6.239	1377	1389	1403	rVB2	48200	131011	1.74%	0.103%
19	6.352	1403	1424	1455	rBV	560246	1756814	23.37%	1.380%
20	6.500	1456	1470	1482	rBV3	64189	179111	2.38%	0.141%
21	6.606	1483	1503	1520	rBV	1399617	4353162	57.92%	3.419%
22	7.181	1661	1682	1702	rVB9	39798	161599	2.15%	0.127%
23	7.307	1702	1721	1726	rBV2	108574	255607	3.40%	0.201%
24	7.371	1728	1741	1761	rVB	912655	2364639	31.46%	1.857%
25	7.574	1781	1804	1810	rBV3	285024	781700	10.40%	0.614%
26	7.644	1812	1826	1840	rVV7	916059	2823421	37.57%	2.218%
27	7.731	1842	1853	1875	rVB6	381355	1110522	14.78%	0.872%
28	7.860	1876	1893	1907	rBV2	485995	1202327	16.00%	0.944%
29	8.024	1925	1944	1965	rBV3	1746006	4442983	59.11%	3.490%
30	8.153	1966	1984	1996	rBV7	516877	1615953	21.50%	1.269%
31	8.288	2019	2026	2047	rVB3	207682	507388	6.75%	0.399%
32	8.397	2047	2060	2088	rVB4	142193	422085	5.62%	0.332%
33	8.570	2095	2114	2136	rBV3	1465789	3991207	53.10%	3.135%
34	8.664	2137	2143	2155	rVB	162119	303845	4.04%	0.239%

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ALS Vial : 10 Sample Multiplier: 1

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ClientSampleId : MW-3

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

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Title : SW846 8260

35	8.734	2155	2165	2176	rVB	153714	289131	3.85%	0.227%
36	8.808	2176	2188	2195	rBV	410335	843227	11.22%	0.662%
37	9.037	2241	2259	2262	rBV2	393210	731589	9.73%	0.575%
38	9.136	2283	2290	2304	rVB4	72608	153052	2.04%	0.120%
39	9.262	2318	2329	2342	rVB	170733	346935	4.62%	0.272%
40	9.355	2343	2358	2369	rBV7	56210	168924	2.25%	0.133%
41	9.429	2370	2381	2384	rBV2	144925	262916	3.50%	0.206%
42	9.509	2398	2406	2417	rVB2	180341	325106	4.33%	0.255%
43	9.606	2424	2436	2445	rBV2	890064	1770741	23.56%	1.391%
44	9.860	2497	2515	2521	rBV8	106177	333500	4.44%	0.262%
45	9.963	2536	2547	2557	rBV	471848	853620	11.36%	0.670%
46	10.014	2558	2563	2572	rVB3	115816	191940	2.55%	0.151%
47	10.085	2573	2585	2597	rBV	1375879	2716178	36.14%	2.133%
48	10.149	2598	2605	2616	rVB	353947	591485	7.87%	0.465%
49	10.287	2631	2648	2662	rBV5	277410	734022	9.77%	0.577%
50	10.381	2673	2677	2685	rVB3	81844	108348	1.44%	0.085%
51	10.516	2709	2719	2736	rVB4	579263	1232656	16.40%	0.968%
52	10.789	2790	2804	2810	rBV3	132132	288940	3.84%	0.227%
53	10.831	2812	2817	2823	rVV3	99394	142360	1.89%	0.112%
54	10.873	2824	2830	2836	rVV5	109848	173334	2.31%	0.136%
55	10.914	2837	2843	2853	rVB2	192777	298814	3.98%	0.235%
56	11.114	2898	2905	2915	rVB3	97097	170985	2.27%	0.134%
57	11.217	2925	2937	2954	rVB10	70079	228381	3.04%	0.179%
58	11.326	2955	2971	2980	rBV9	82854	230317	3.06%	0.181%
59	11.410	2985	2997	3009	rBV	1170737	2120665	28.22%	1.666%
60	11.512	3019	3029	3047	rVB	1141298	2085902	27.75%	1.638%
61	11.622	3049	3063	3084	rVB	1051922	2199688	29.27%	1.728%
62	11.757	3098	3105	3130	rVB	52983	170611	2.27%	0.134%
63	11.866	3130	3139	3148	rBV3	50287	105538	1.40%	0.083%
64	11.950	3158	3165	3183	rVB	137337	255374	3.40%	0.201%
65	12.252	3247	3259	3273	rBV	2642073	4310943	57.36%	3.386%
66	12.406	3295	3307	3323	rVB	898165	1562413	20.79%	1.227%
67	12.596	3354	3366	3382	rBV	1871544	3048186	40.56%	2.394%
68	12.738	3402	3410	3424	rVB	365615	595821	7.93%	0.468%
69	12.895	3448	3459	3474	rBV	818497	1317780	17.53%	1.035%
70	13.046	3499	3506	3515	rVB	75965	118546	1.58%	0.093%
71	13.175	3532	3546	3558	rBV3	205523	478843	6.37%	0.376%

LSC Area Percent Report

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ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

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

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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Title : SW846 8260

72	13.348	3589	3600	3620	rVB2	1024219	2002954	26.65%	1.573%
73	13.448	3624	3631	3640	rVB	85919	131958	1.76%	0.104%
74	13.519	3641	3653	3656	rBV	1562540	2266042	30.15%	1.780%
75	13.551	3656	3663	3672	rVV	3513508	5906563	78.59%	4.639%
76	13.599	3672	3678	3693	rVB2	491827	967639	12.87%	0.760%
77	13.680	3693	3703	3713	rVB	373681	567297	7.55%	0.446%
78	13.747	3714	3724	3734	rVB	261374	416977	5.55%	0.327%
79	13.895	3757	3770	3780	rBV	3151687	4898399	65.17%	3.847%
80	13.953	3780	3788	3794	rVV	400659	621761	8.27%	0.488%
81	14.001	3794	3803	3816	rVB2	2046671	3314007	44.09%	2.603%
82	14.139	3836	3846	3854	rBV2	86589	136633	1.82%	0.107%
83	14.217	3855	3870	3888	rBV	2114562	3347213	44.53%	2.629%
84	14.300	3888	3896	3903	rBV2	214585	334817	4.45%	0.263%
85	14.374	3912	3919	3927	rBV2	169548	228484	3.04%	0.179%
86	14.438	3932	3939	3944	rVV	110095	164740	2.19%	0.129%
87	14.483	3944	3953	3963	rVV2	1045246	1711703	22.77%	1.344%
88	14.535	3963	3969	3975	rVB	128913	172459	2.29%	0.135%
89	14.596	3975	3988	4000	rBV3	1472252	3118127	41.49%	2.449%
90	14.664	4000	4009	4020	rVB4	228781	423895	5.64%	0.333%
91	14.860	4053	4070	4079	rBV3	574546	1178656	15.68%	0.926%
92	14.966	4091	4103	4119	rVB2	523723	1105565	14.71%	0.868%
93	15.133	4146	4155	4166	rVB	348168	575138	7.65%	0.452%
94	15.233	4178	4186	4193	rBV2	112186	193273	2.57%	0.152%
95	15.393	4227	4236	4247	rBV	191938	306758	4.08%	0.241%
96	15.538	4273	4281	4305	rVB	122418	282958	3.76%	0.222%
97	15.644	4306	4314	4323	rBV	98543	161956	2.15%	0.127%
98	15.708	4328	4334	4347	rVB2	99555	175910	2.34%	0.138%
99	16.024	4421	4432	4461	rVB3	389895	889123	11.83%	0.698%
100	16.217	4482	4492	4506	rVB	185875	362932	4.83%	0.285%

Sum of corrected areas: 127321884

LSC Report - Integrated Chromatogram

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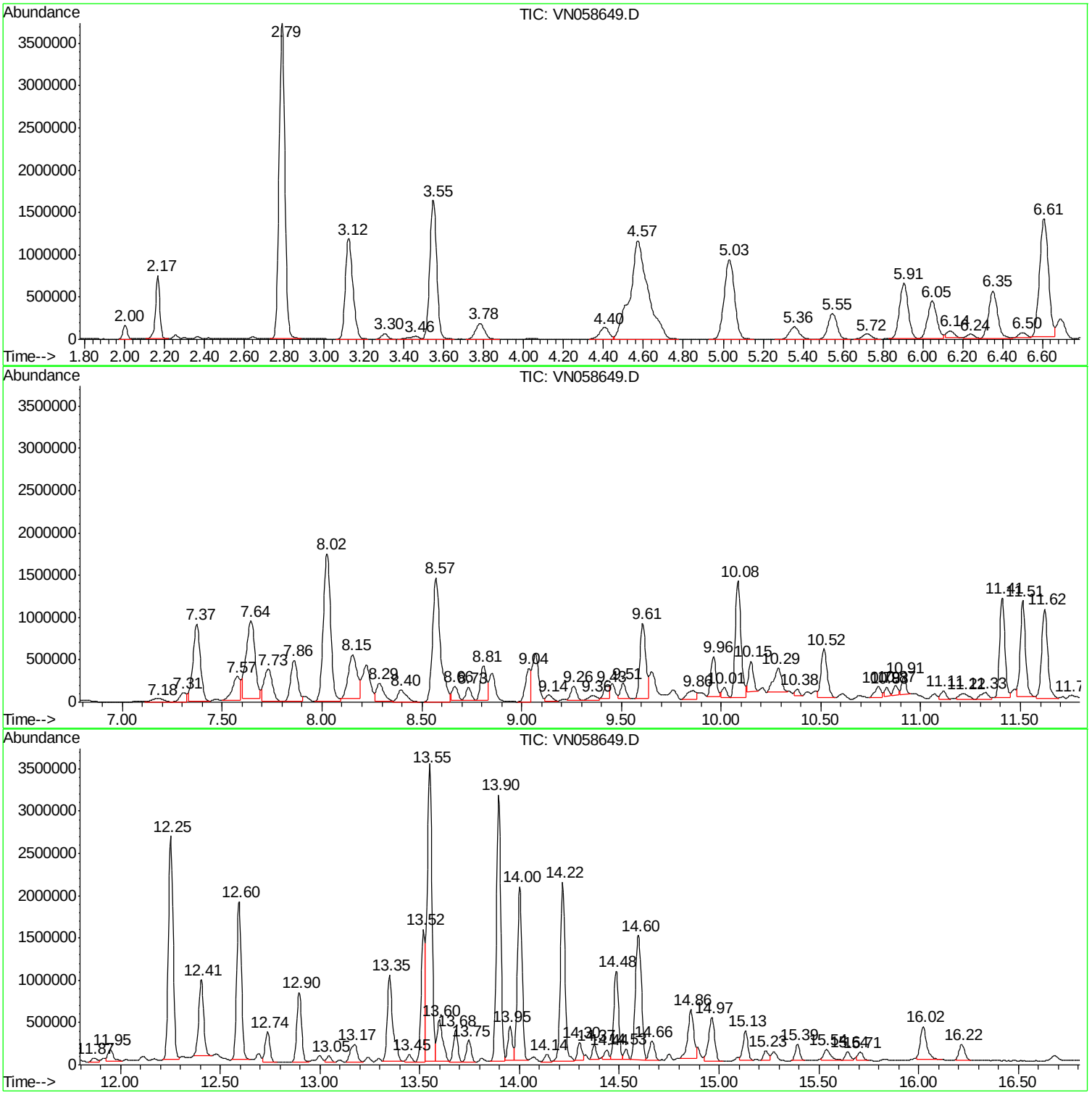
Sample : K5184-08
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260

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



Library Search Compound Report

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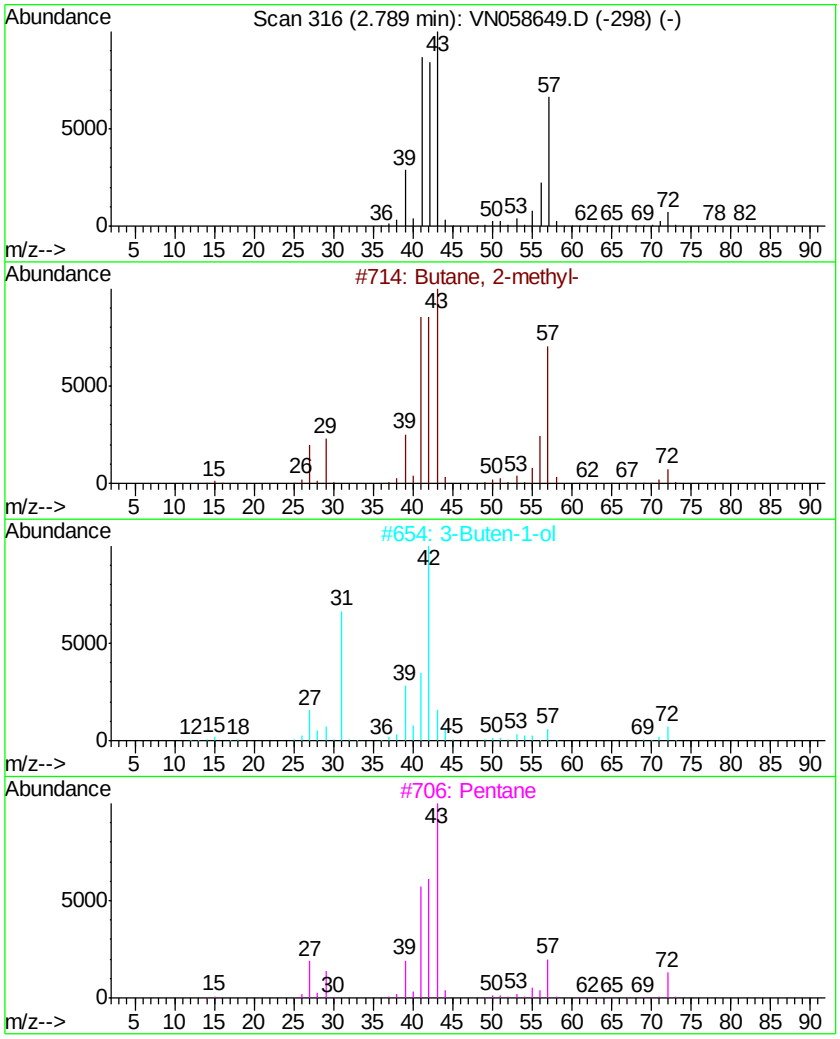
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

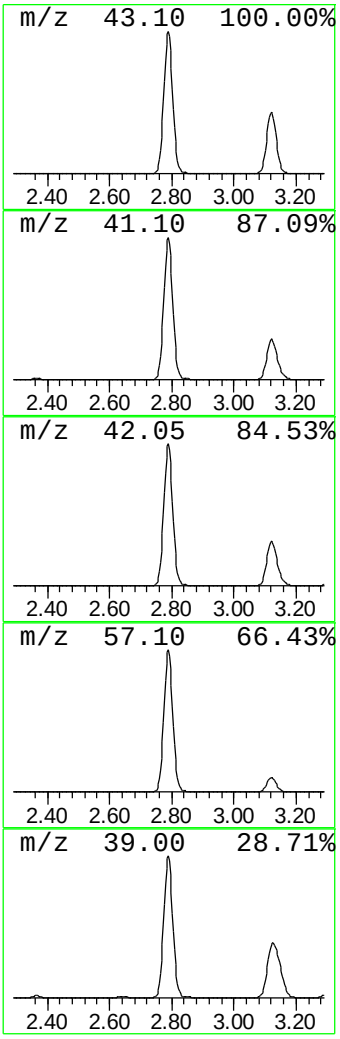
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.79	133.10 ug/l	7516050	Pentafluorobenzene	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	94
2	3-Buten-1-ol	72	C4H8O	000627-27-0	47
3	Pentane	72	C5H12	000109-66-0	45
4	1-Butene	56	C4H8	000106-98-9	9
5	Methane, isocyanato-	57	C2H3NO	000624-83-9	9





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ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260

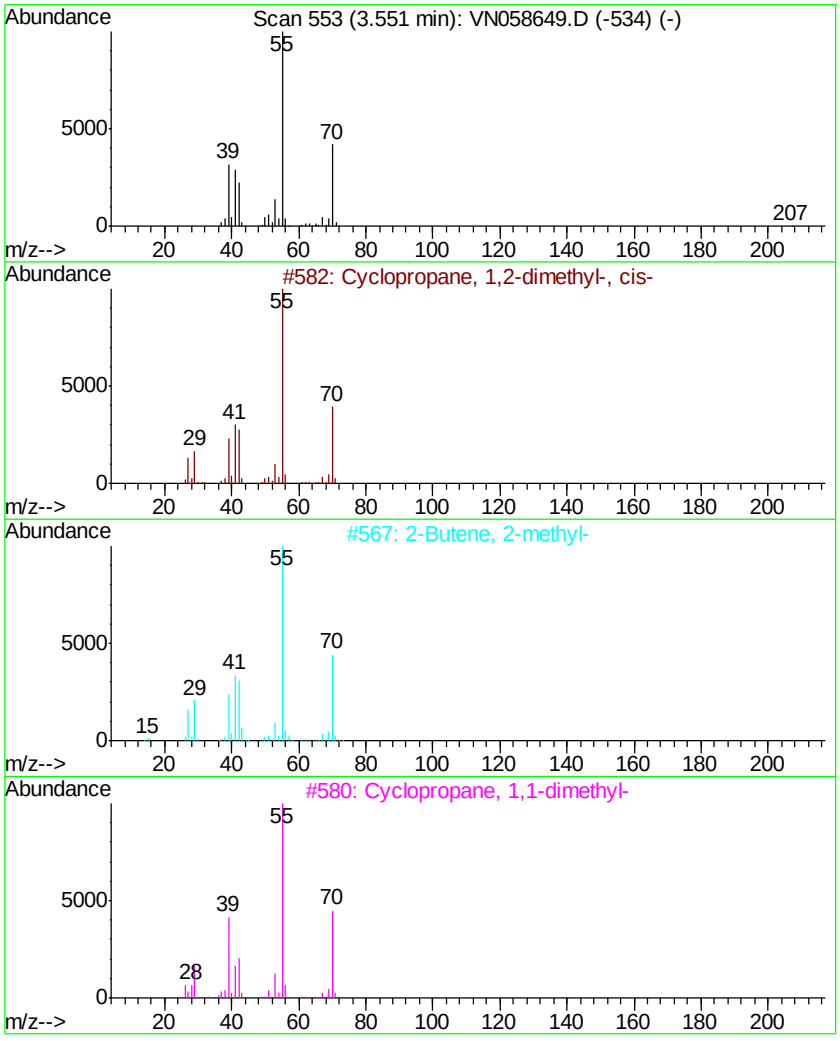
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

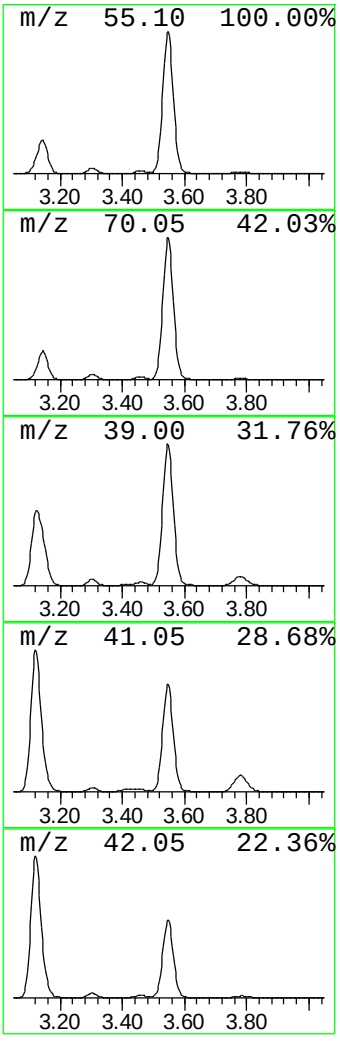
****

Peak Number 2 Cyclopropane, 1,2-dimethyl-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.55	69.85 ug/l	3944450	Pentafluorobenzene	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	91
2	2-Butene, 2-methyl-	70	C5H10	000513-35-9	90
3	Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	80
4	Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	76
5	2-Pentene, (Z)-	70	C5H10	000627-20-3	68





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Quant Title : SW846 8260

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

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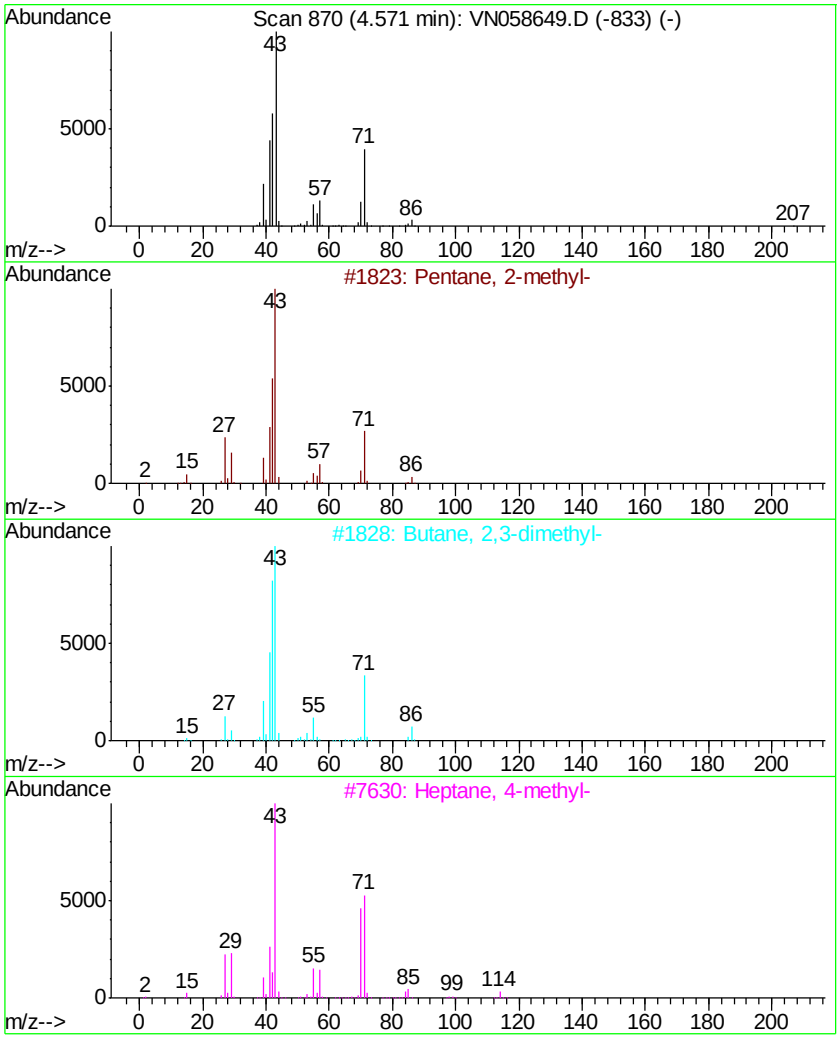
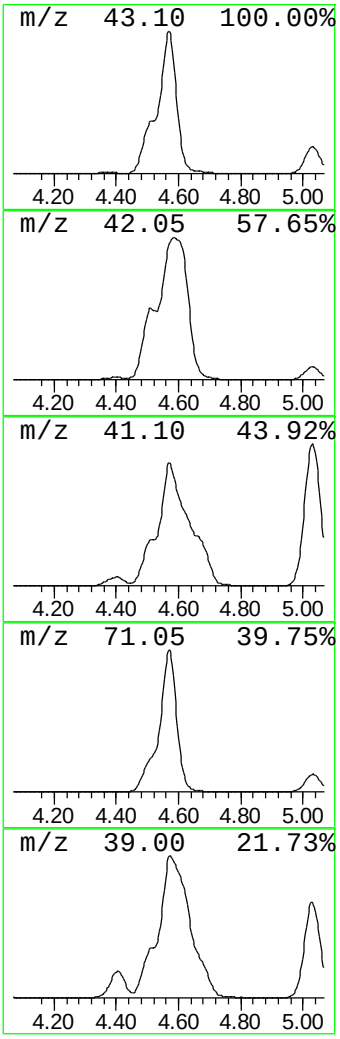
Peak Number 3 Pentane, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.57	130.67 ug/l	7378930	Pentafluorobenzene	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	72
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	28
4	C2H5C(CH3)2COCH3	114	C7H14O	020669-04-9	25
5	Pentane	72	C5H12	000109-66-0	25





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Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260

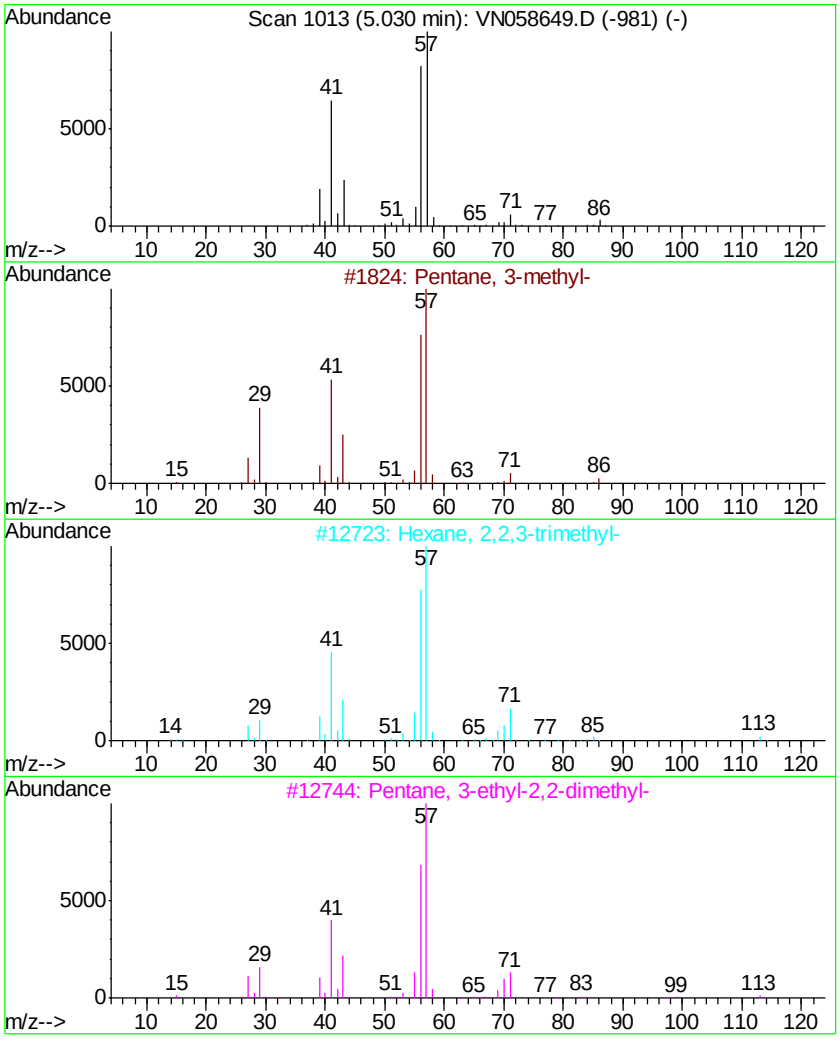
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

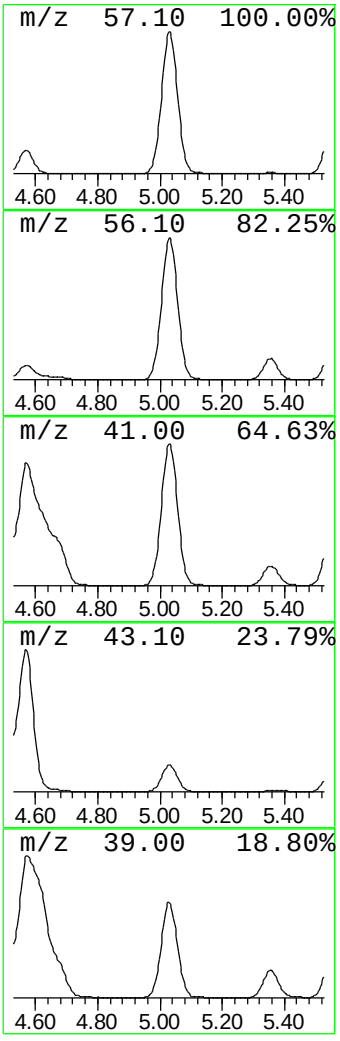
****

Peak Number 4 Pentane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.03	60.37 ug/l	3408810	Pentafluorobenzene	7.65

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	83
3	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
4	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45
5	Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	40





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ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260

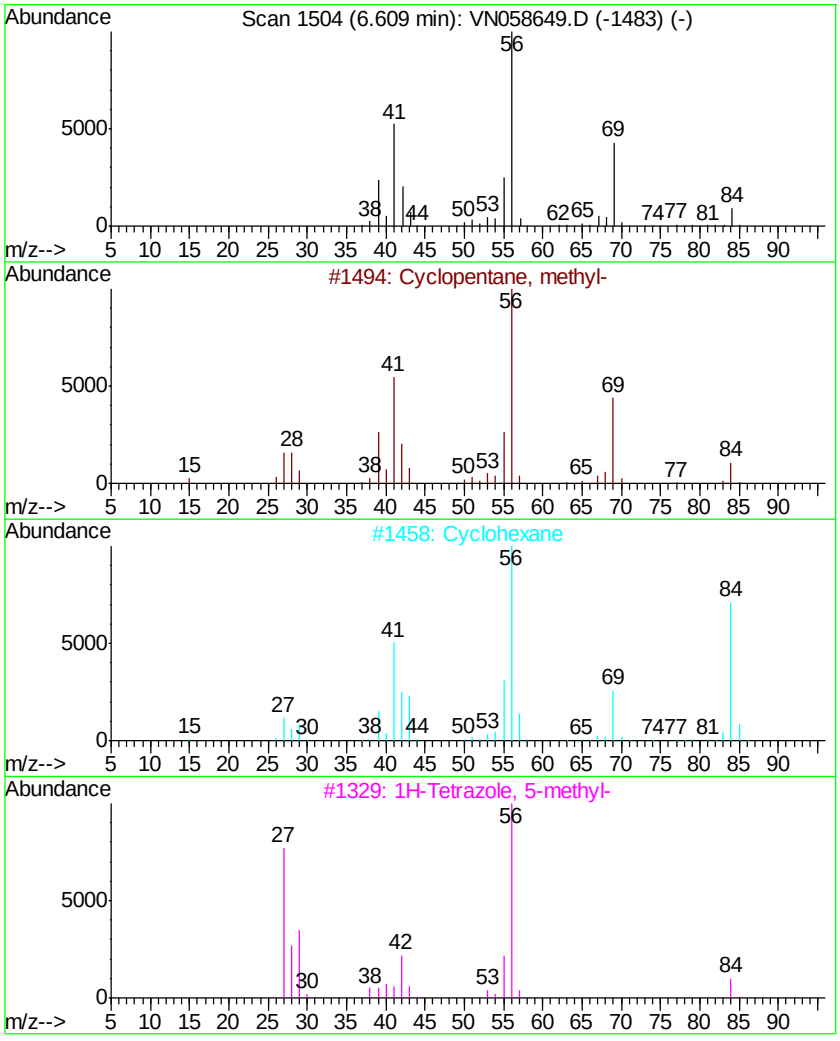
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

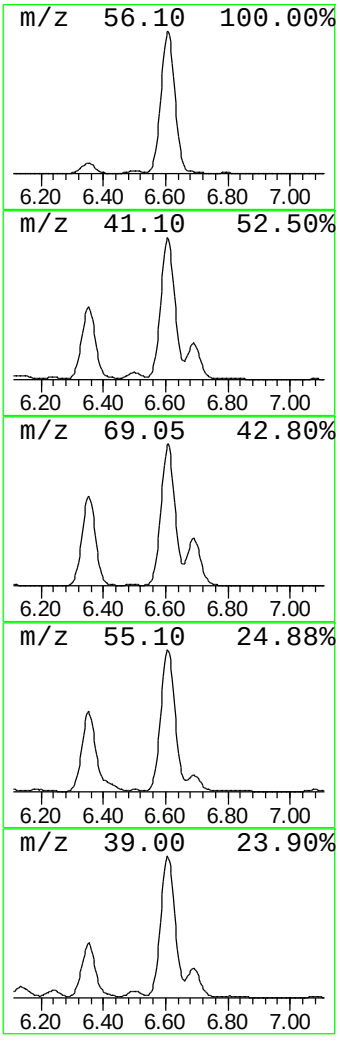
****

Peak Number 5 Cyclopentane, methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	77.09 ug/l	4353160	Pentafluorobenzene	7.65

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	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	56





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN101019\
Data File : VN058649.D
Acq On : 9 Oct 2019 18:48
Operator : JC/SP
Sample : K5184-08
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260

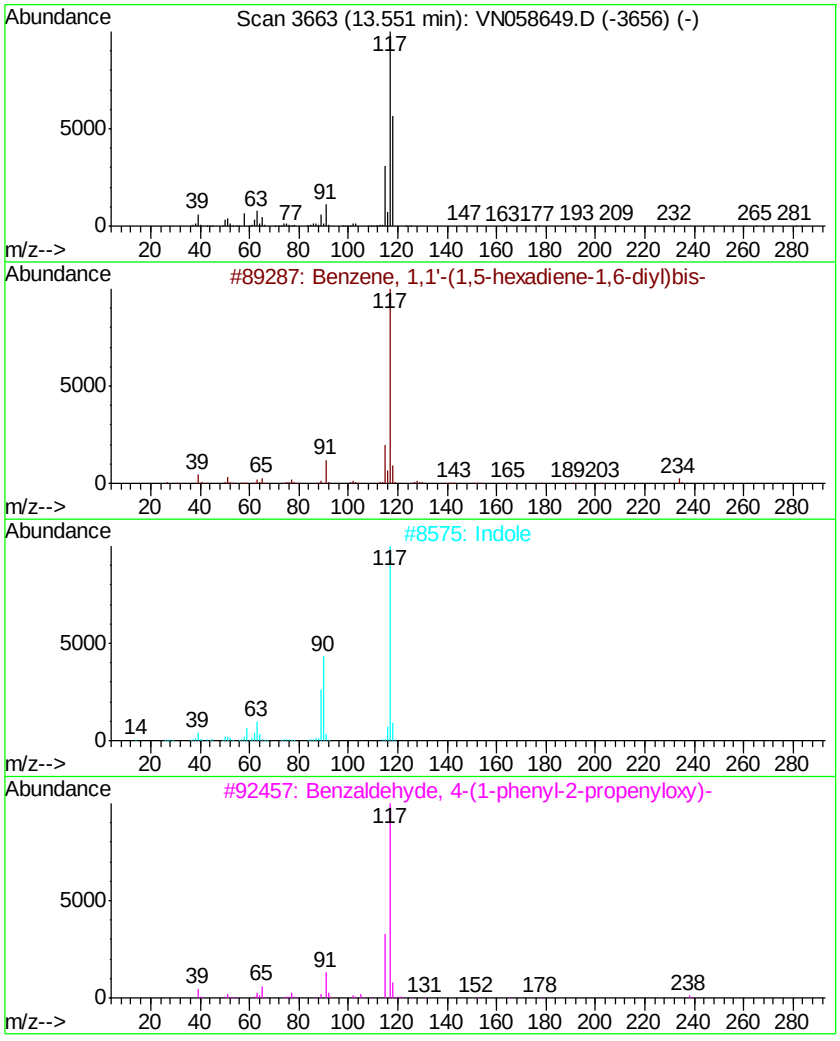
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

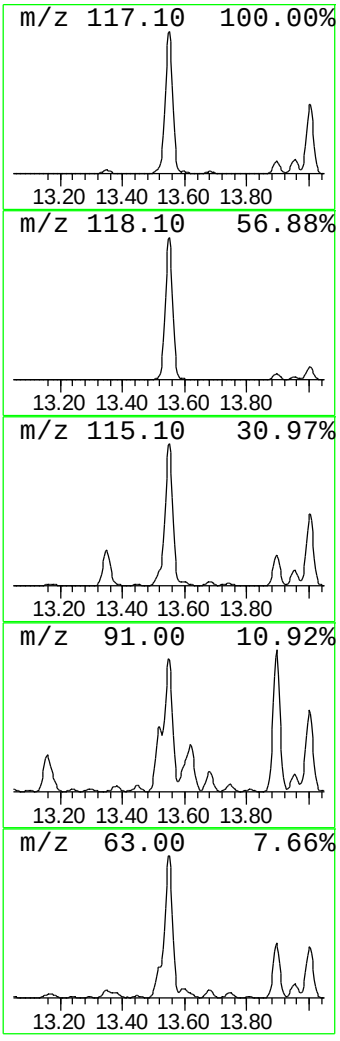
****

Peak Number 6 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.55	147.45 ug/l	5906560	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	53
2	Indole	117	C8H7N	000120-72-9	46
3	Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	42
4	trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	40
5	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	32





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Data File : VN058649.D
Acq On : 9 Oct 2019 18:48
Operator : JC/SP
Sample : K5184-08
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260

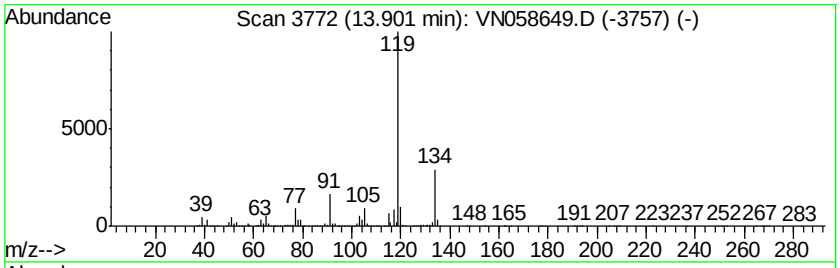
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

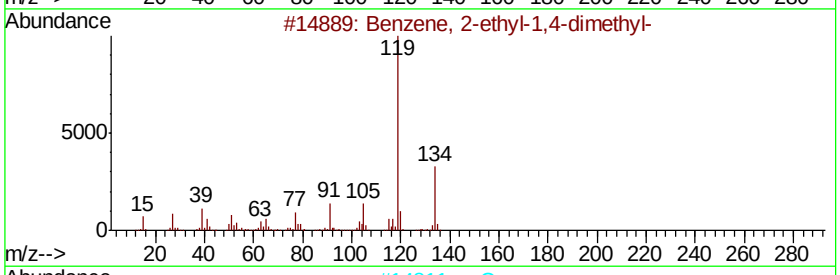
****

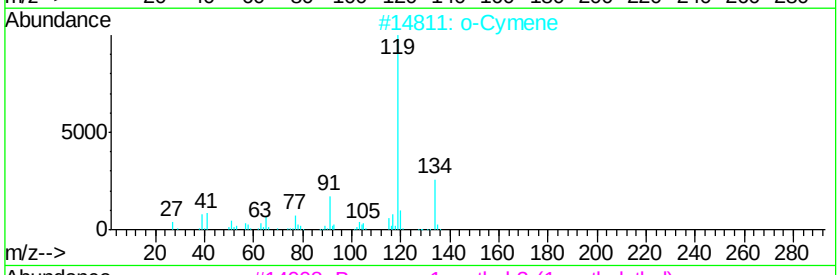
Peak Number 7 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 4

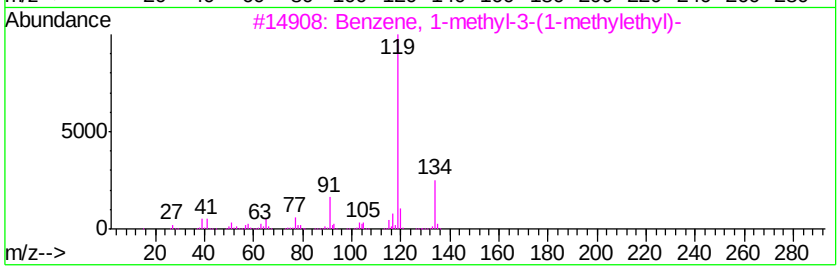
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	122.28 ug/l	4898400	1,4-Dichlorobenzene-d4	13.35

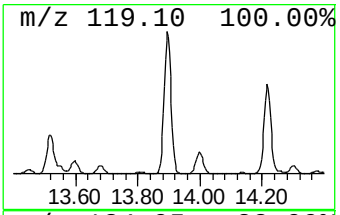
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94

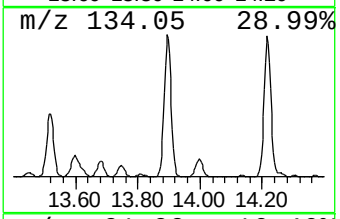


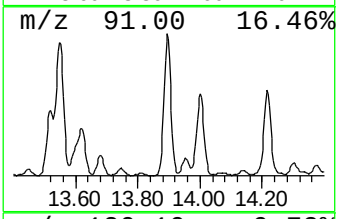


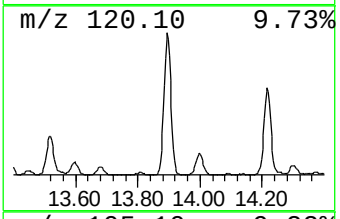


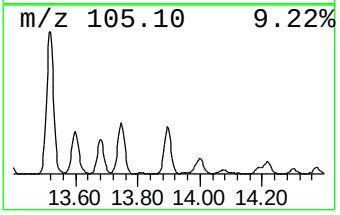












Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN101019\
Data File : VN058649.D
Acq On : 9 Oct 2019 18:48
Operator : JC/SP
Sample : K5184-08
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260

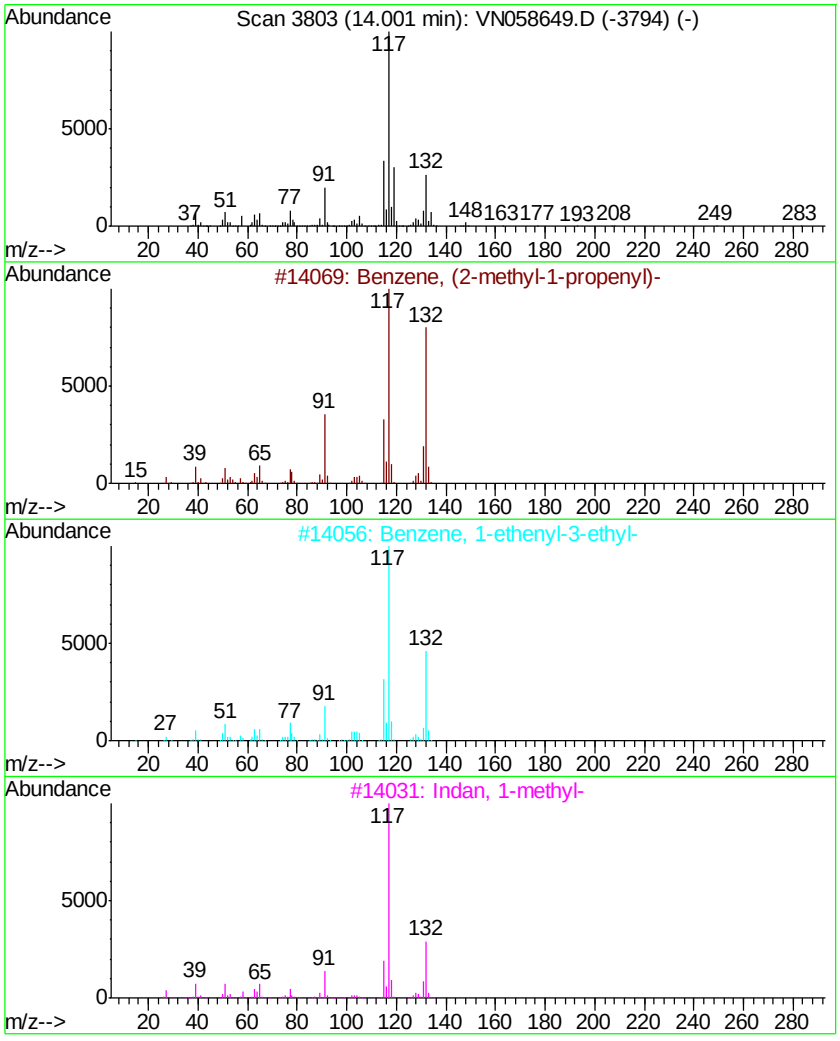
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

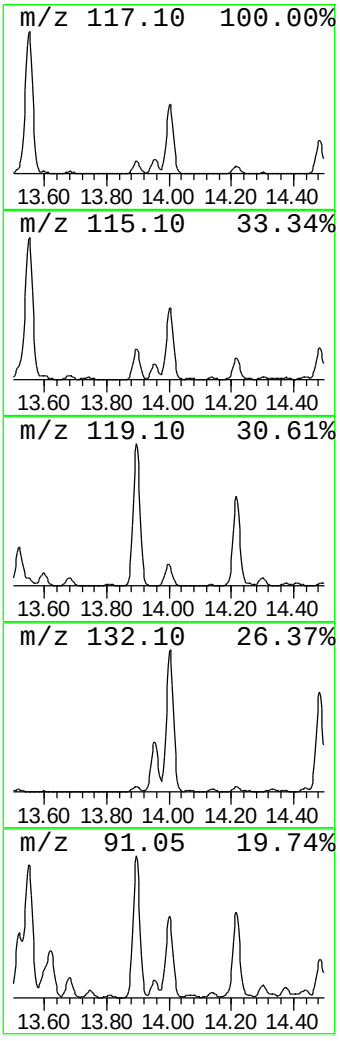
****

Peak Number 8 Benzene, (2-methyl-1-propen... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.00	82.73 ug/l	3314010	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	81
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
3	Indan, 1-methyl-	132	C10H12	000767-58-8	76
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	74





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN101019\
Data File : VN058649.D
Acq On : 9 Oct 2019 18:48
Operator : JC/SP
Sample : K5184-08
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260

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

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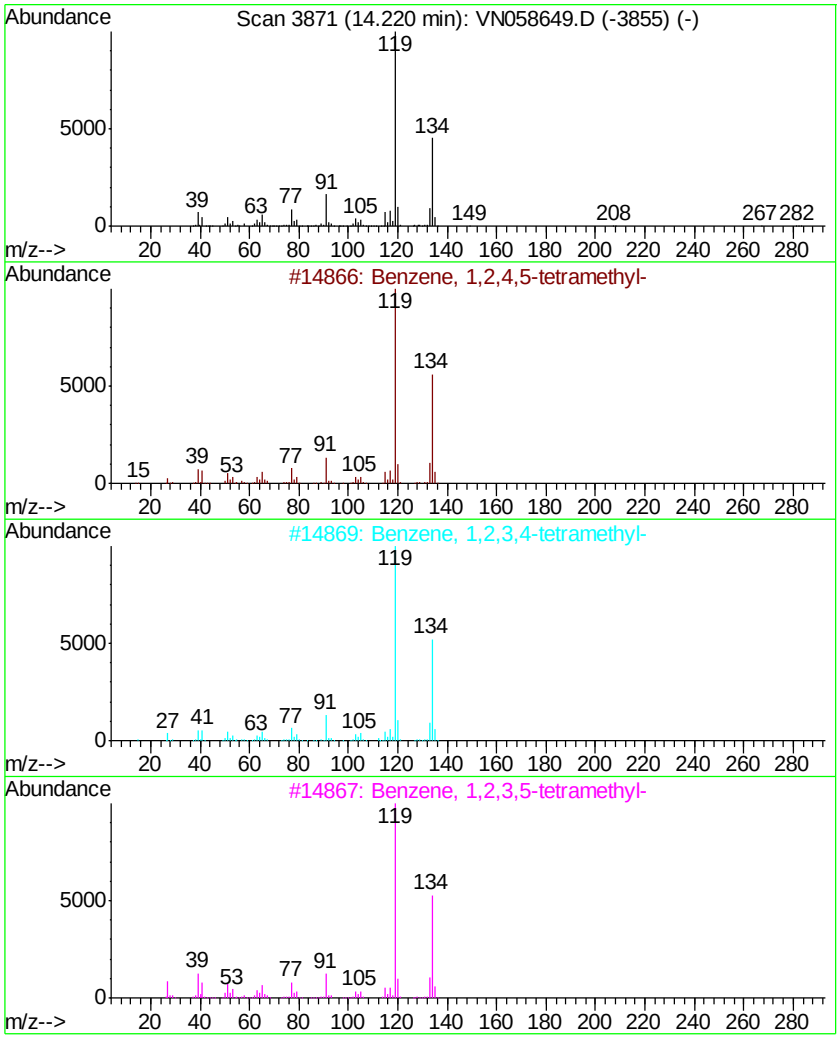
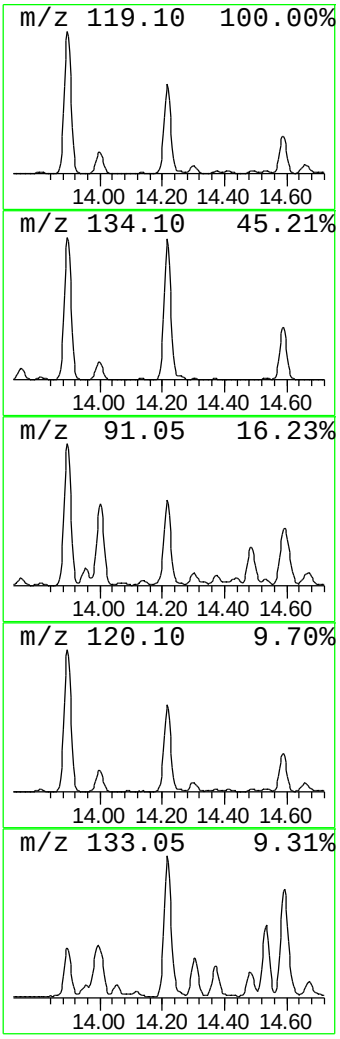
Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.22	83.56 ug/l	3347210	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94





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Data File : VN058649.D
Acq On : 9 Oct 2019 18:48
Operator : JC/SP
Sample : K5184-08
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260

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

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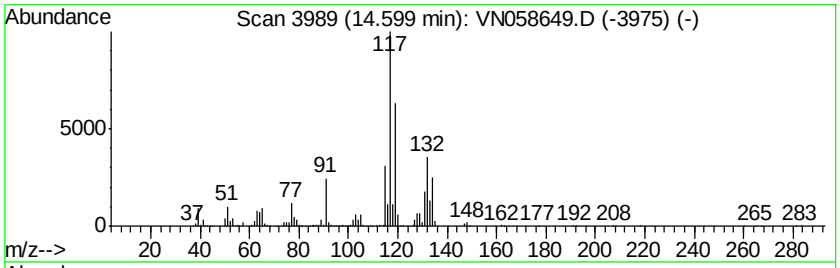
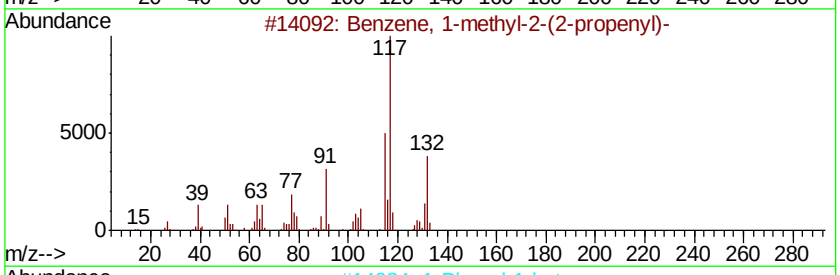
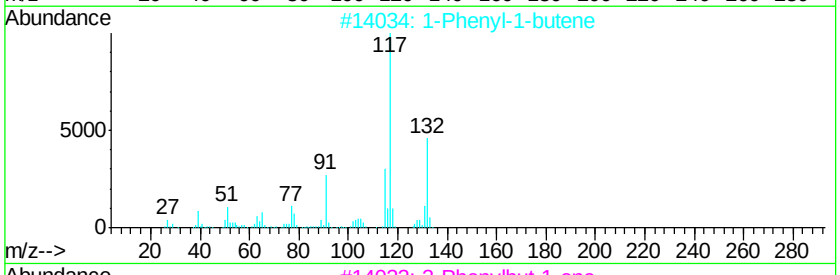
Peak Number 10 Benzene, 1-methyl-2-(2-prop... Concentration Rank 7

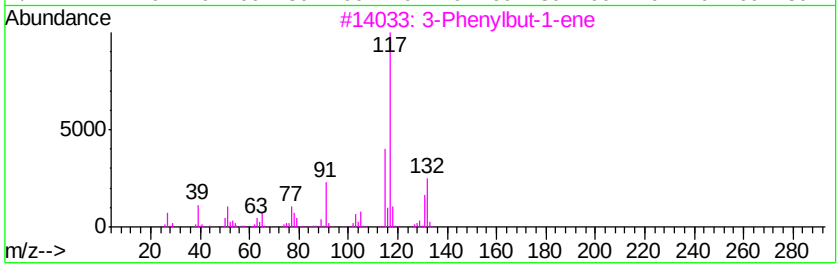
R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	77.84 ug/l	3118130	1,4-Dichlorobenzene-d4	13.35

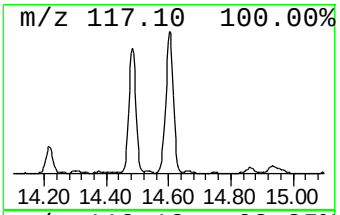
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	86
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	86
4	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	70
5	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70

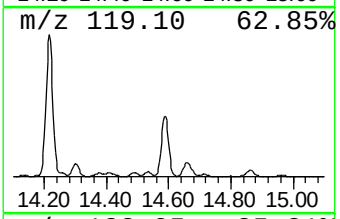


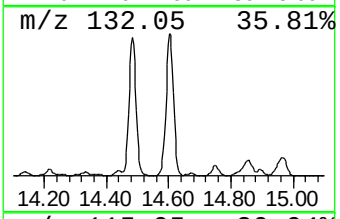


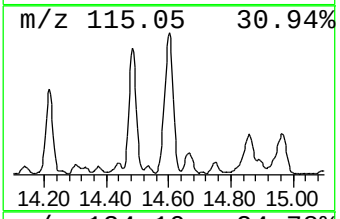


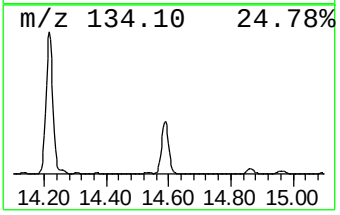












Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN101019\
Data File : VN058649.D
Acq On : 9 Oct 2019 18:48
Operator : JC/SP
Sample : K5184-08
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methyl-	2.79	133.1	ug/l	7516050	1	7.65	2823420	50.0
Cyclopropane, 1,2...	3.55	69.8	ug/l	3944450	1	7.65	2823420	50.0
Pentane, 2-methyl-	4.57	130.7	ug/l	7378930	1	7.65	2823420	50.0
Pentane, 3-methyl-	5.03	60.4	ug/l	3408810	1	7.65	2823420	50.0
Cyclopentane, met...	6.61	77.1	ug/l	4353160	1	7.65	2823420	50.0
Benzene, 1,1'-(1,...	13.55	147.4	ug/l	5906560	4	13.35	2002950	50.0
Benzene, 2-ethyl-...	13.90	122.3	ug/l	4898400	4	13.35	2002950	50.0
Benzene, (2-methy...	14.00	82.7	ug/l	3314010	4	13.35	2002950	50.0
Benzene, 1,2,4,5-...	14.22	83.6	ug/l	3347210	4	13.35	2002950	50.0
Benzene, 1-methyl...	14.60	77.8	ug/l	3118130	4	13.35	2002950	50.0