LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.911	3	4	16	rVB	114612	132561	2.36%	0.327%
2	3.193	48	52	59	rBV	171639	218408	3.89%	0.539%
3	3.887	167	170	175	rVB	12300	16579	0.30%	0.041%
4	3.964	179	183	187	rVV	41688	56456	1.01%	0.139%
5	4.005	188	190	193	rVB2	8132	8223	0.15%	0.020%
6	4.311	238	242	251	rVB	48489	79055	1.41%	0.195%
7	4.987	353	357	362	rVB2	9758	14277	0.25%	0.035%
8	5.681	467	475	483	rBV8	3834	9717	0.17%	0.024%
9	5.928	511	517	523	rVB	9269	14684	0.26%	0.036%
10	6.346	582	588	596	rVB	89045	136410	2.43%	0.336%
11	6.599	626	631	635	rBV7	13005	26224	0.47%	0.065%
12	6.646	635	639	646	rVB	31157	50806	0.90%	0.125%
13	6.828	660	670	683	rVB3	18922	50291	0.90%	0.124%
14	6.975	688	695	705	rBV	61046	106546	1.90%	0.263%
15	7.099	710	716	721	rBV9	7498	13901	0.25%	0.034%
16	7.175	723	729	737	rBV	65015	115994	2.07%	0.286%
17	7.381	758	764	768	rBV9	5913	10307	0.18%	0.025%
18	7.546	787	792	796	rVB8	5410	8246	0.15%	0.020%
19	7.634	798	807	816	rBV	1030011	1662725	29.61%	4.100%
20	7.705	816	819	826	rVV6	14926	33796	0.60%	0.083%
21	7.781	826	832	842	rVB	60134	98758	1.76%	0.244%
22	7.869	842	847	852	rBV	37596	58755	1.05%	0.145%
23	7.946	854	860	865	rBV2	28961	46222	0.82%	0.114%
24	8.293	912	919	922	rBV9	3653	7863	0.14%	0.019%
25	8.352	924	929	936	rVV	49548	90329	1.61%	0.223%
26	8.416	936	940	944	rVB6	6885	11640	0.21%	0.029%
27	8.516	952	957	962	rBV9	4526	10978	0.20%	0.027%
28	8.611	966	973	989	rVV	1014859	1672403	29.78%	4.124%
29	8.734	989	994	996	rVV	24180	40592	0.72%	0.100%
30	8.781	996	1002	1011	rVV	64976	146387	2.61%	0.361%
31	8.869	1011	1017	1024	rVB	216863	339697	6.05%	0.838%
32	8.981	1031	1036	1042	rVB8	9723	18515	0.33%	0.046%
33	9.140	1057	1063	1068	rBV7	17780	31663	0.56%	0.078%
34	9.187	1068	1071	1076	rVV5	6340	9189	0.16%	0.023%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Title : SVOA CALIBRATION

35	9.252	1076	1082	1089	rVV3	20969	43305	0.77%	0.107%
36	9.340	1089	1097	1106	rVV	592700	968366	17.25%	2.388%
37	9.440	1110	1114	1118	rVV7	8525	14959	0.27%	0.037%
38	9.505	1119	1125	1129	rVV3	36378	73889	1.32%	0.182%
39	9.546	1129	1132	1138	rVB3	20705	33775	0.60%	0.083%
40	9.628	1141	1146	1149	rVV4	13234	25402	0.45%	0.063%
41	9.675	1151	1154	1160	rVV2	22963	38543	0.69%	0.095%
42	9.763	1160	1169	1172	rVV	259397	488551	8.70%	1.205%
43	9.828	1175	1180	1189	rVB	749021	1231668	21.94%	3.037%
44	9.963	1195	1203	1209	rVB	79302	150071	2.67%	0.370%
45	10.046	1209	1217	1219	rBV	120294	205905	3.67%	0.508%
46	10.187	1234	1241	1251	rBV	515383	856968	15.26%	2.113%
47	10.281	1251	1257	1266	rVV	141002	233214	4.15%	0.575%
48	10.405	1271	1278	1284	rVV	1412663	2401814	42.77%	5.923%
49	10.475	1284	1290	1297	rVV	2096175	3265252	58.15%	8.052%
50	10.546	1297	1302	1314	rVB2	125813	281658	5.02%	0.695%
51	10.987	1372	1377	1380	rBV7	7724	15538	0.28%	0.038%
52	11.105	1395	1397	1403	rVB7	11982	15991	0.28%	0.039%
53	11.246	1409	1421	1432	rBV	181336	363008	6.46%	0.895%
54	11.387	1439	1445	1452	rVB10	12817	27369	0.49%	0.067%
55	11.457	1453	1457	1460	rBV6	5222	7921	0.14%	0.020%
56	11.610	1479	1483	1489	rVB3	13728	23952	0.43%	0.059%
57	11.769	1503	1510	1515	rBV	25252	45330	0.81%	0.112%
58	11.916	1528	1535	1543	rBV	240085	406883	7.25%	1.003%
59	11.975	1543	1545	1549	rVV5	8974	14412	0.26%	0.036%
60	12.022	1549	1553	1560	rVB	45160	81684	1.45%	0.201%
61	12.152	1567	1575	1585	rBV	128348	216003	3.85%	0.533%
62	12.299	1596	1600	1605	rBV8	7009	13644	0.24%	0.034%
63	12.422	1616	1621	1628	rBV3	19808	39294	0.70%	0.097%
64	12.510	1633	1636	1640	rVB5	7967	11464	0.20%	0.028%
65	12.622	1652	1655	1660	rBV7	5422	8666	0.15%	0.021%
66	12.799	1680	1685	1688	rBV7	7356	13349	0.24%	0.033%
67	12.981	1712	1716	1722	rBV9	9946	18512	0.33%	0.046%
68	13.081	1729	1733	1736	rVB6	6544	8569	0.15%	0.021%
69	13.122	1736	1740	1747	rVB	21844	35198	0.63%	0.087%
70	13.193	1748	1752	1759	rBV8	10928	20225	0.36%	0.050%
71	13.310	1768	1772	1784	rVB8	12942	29824	0.53%	0.074%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Title : SVOA CALIBRATION

72	13.575	1815	1817	1823	rBV7	6535	9549	0.17%	0.024%
73	13.687	1831	1836	1841	rBV	197249	279176	4.97%	0.688%
74	13.734	1841	1844	1851	rVB	45597	61477	1.09%	0.152%
75	13.869	1865	1867	1875	rVB9	8414	14895	0.27%	0.037%
76	13.963	1876	1883	1900	rVV	211006	358257	6.38%	0.884%
77	14.087	1901	1904	1909	rVB6	9309	14540	0.26%	0.036%
78	14.193	1918	1922	1928	rVB2	46506	66004	1.18%	0.163%
79	14.269	1928	1935	1944	rBV2	2313930	3516261	62.62%	8.671%
80	15.028	2060	2064	2068	rBV	30956	48373	0.86%	0.119%
81	15.269	2099	2105	2112	rBV	315641	453826	8.08%	1.119%
82	15.528	2144	2149	2153	rBV7	20901	32093	0.57%	0.079%
83	15.787	2187	2193	2203	rBV	313071	437417	7.79%	1.079%
84	15.957	2217	2222	2228	rBV3	26788	54931	0.98%	0.135%
85	16.122	2246	2250	2254	rBV	15305	22910	0.41%	0.056%
86	16.175	2256	2259	2266	rVB8	9424	13264	0.24%	0.033%
87	16.298	2275	2280	2288	rBV	180948	263865	4.70%	0.651%
88	16.575	2323	2327	2334	rBV2	20875	31511	0.56%	0.078%
89	16.810	2364	2367	2376	rVB2	15703	35192	0.63%	0.087%
90	17.022	2397	2403	2414	rBV	2946919	4315610	76.86%	10.643%
91	17.122	2415	2420	2431	rVB2	347323	519458	9.25%	1.281%
92	18.216	2602	2606	2611	rBV8	13865	21071	0.38%	0.052%
93	18.681	2681	2685	2689	rBV	51080	66848	1.19%	0.165%
94	19.369	2797	2802	2807	rBV	473197	620881	11.06%	1.531%
95	19.422	2808	2811	2817	rVB	42740	56604	1.01%	0.140%
96	21.116	3094	3099	3108	rBV	4268666	5354776	95.37%	13.205%
97	22.686	3363	3366	3374	rVB	211844	308721	5.50%	0.761%
98	23.180	3446	3450	3456	rVB	309424	537261	9.57%	1.325%
99	23.251	3459	3462	3467	rVV	228397	341529	6.08%	0.842%
100	23.322	3468	3474	3487	rVB2	3004846	5615016	100.00%	13.847%

Sum of corrected areas: 40549689

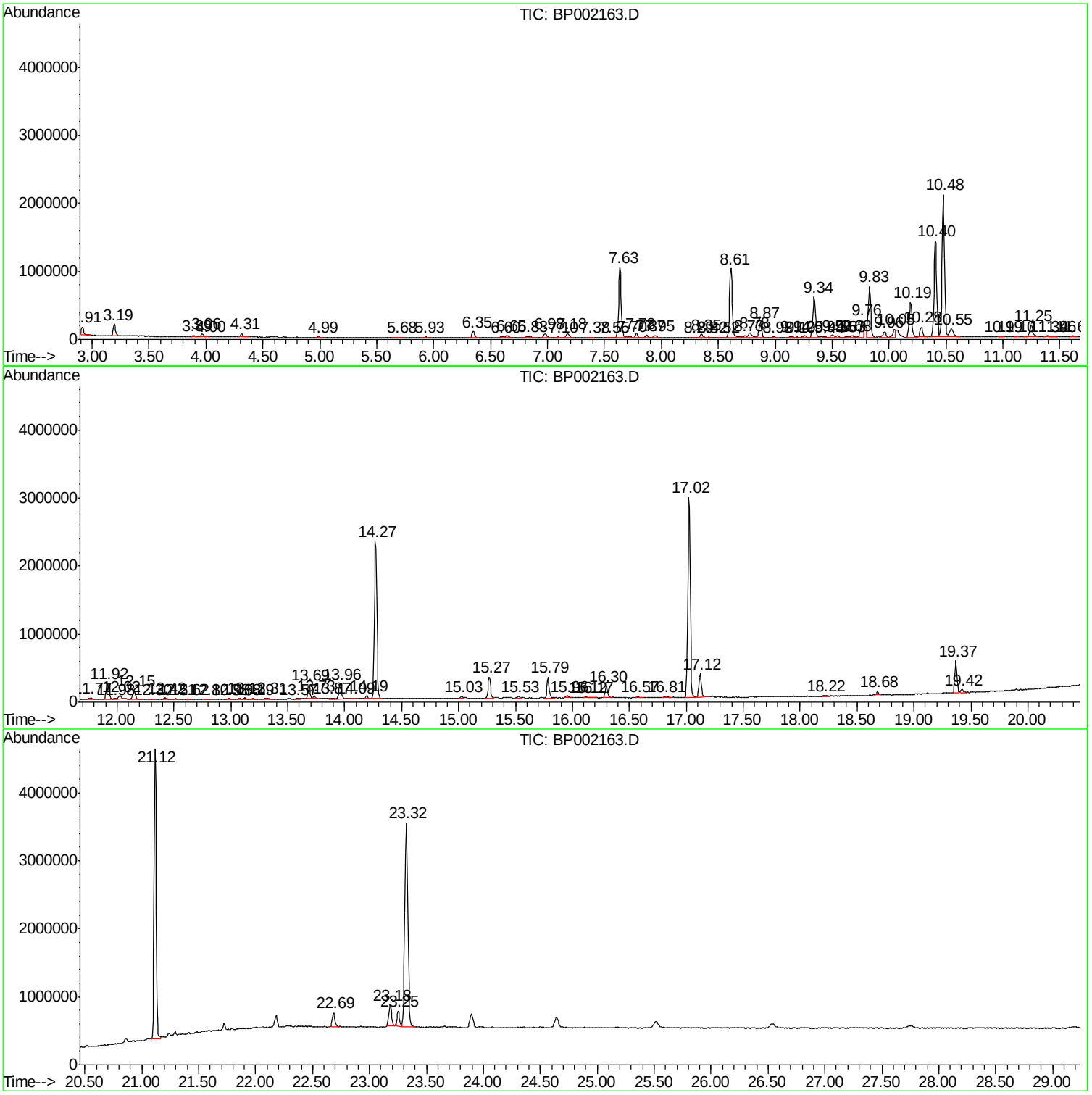
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

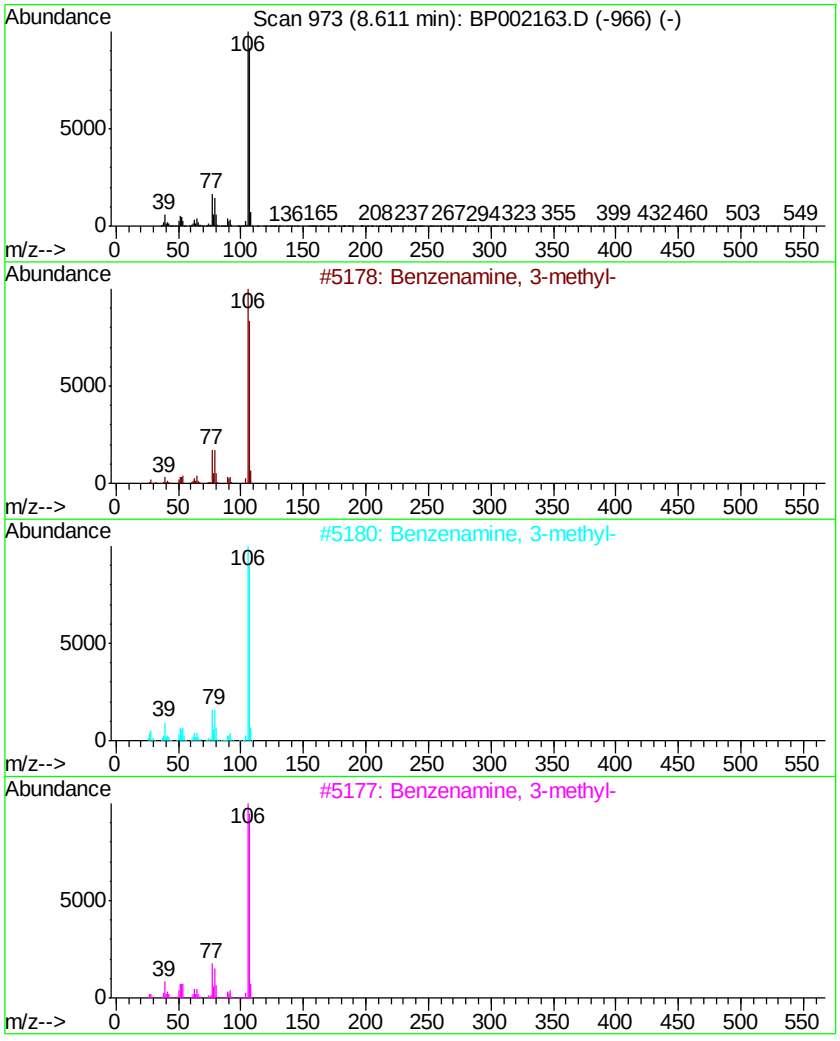
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

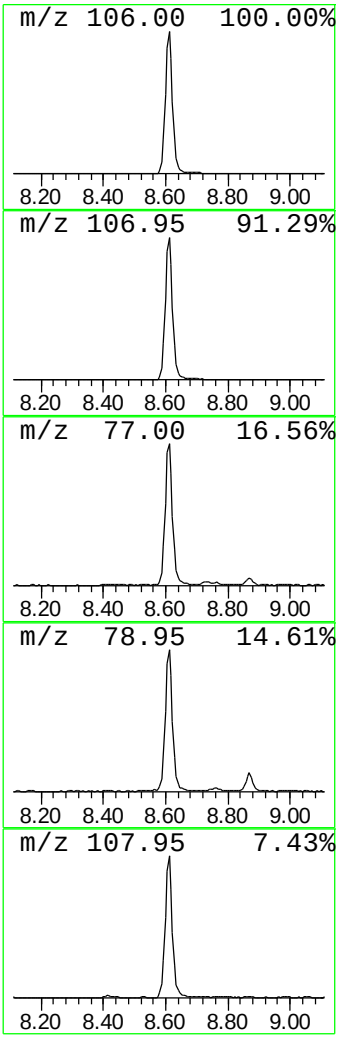
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.61	20.12 ng/ul	1672400	1,4-Dichlorobenzene-d4	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
2	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
3	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
4	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	95
5	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	94





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

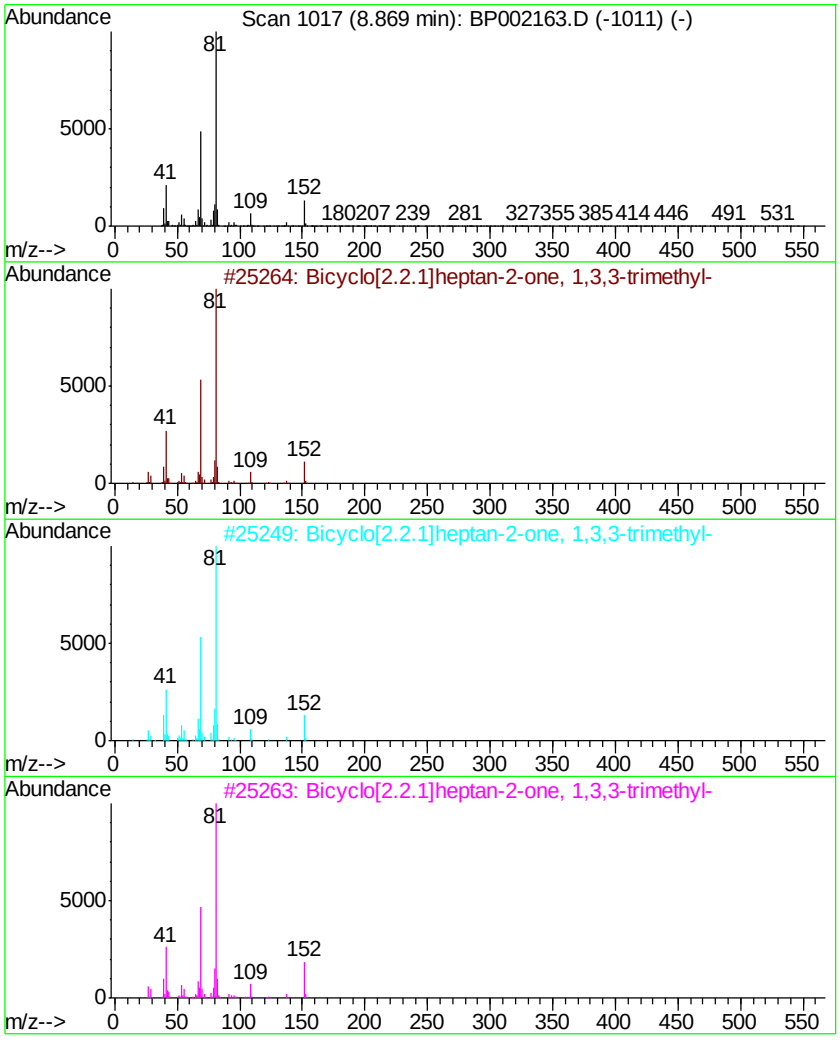
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

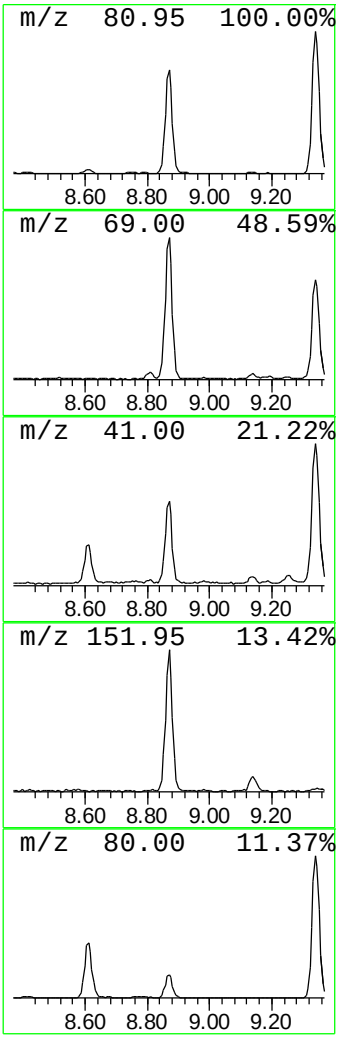
****

Peak Number 2 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.87	4.09 ng/ul	339697	1,4-Dichlorobenzene-d4	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	95
2	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	95
3	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	95
4	D-Fenchone	152	C10H16O	004695-62-9	94
5	L-Fenchone	152	C10H16O	007787-20-4	94





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

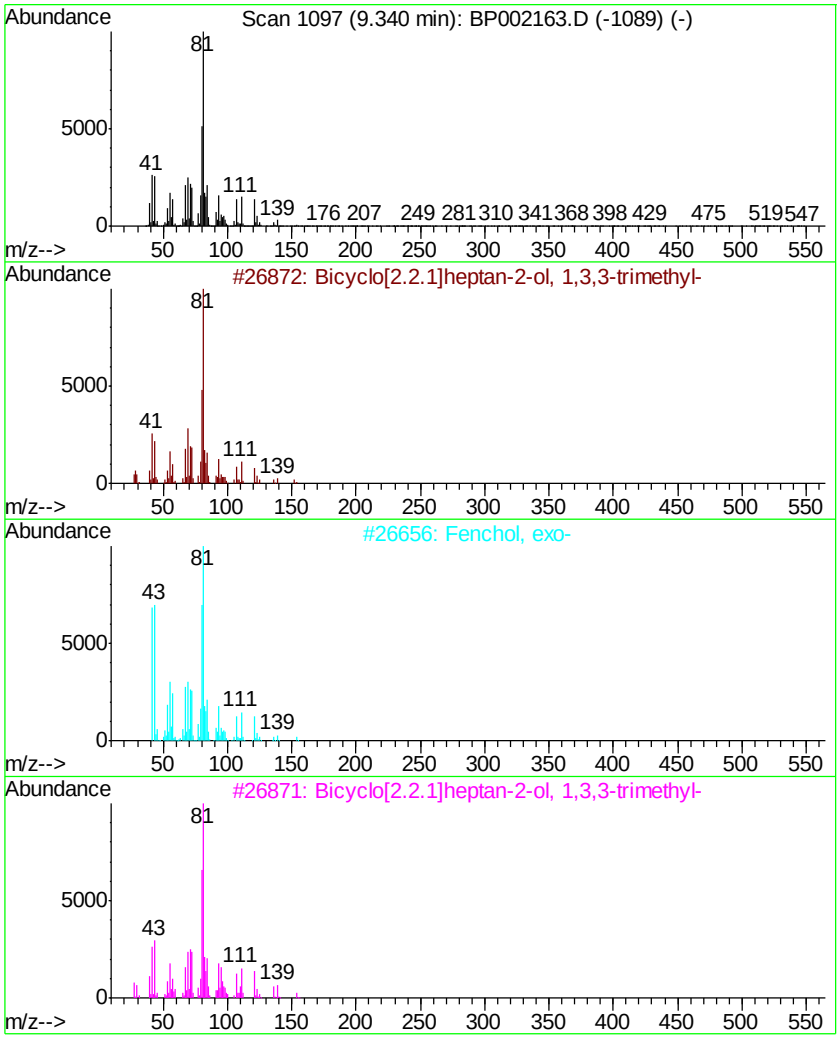
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

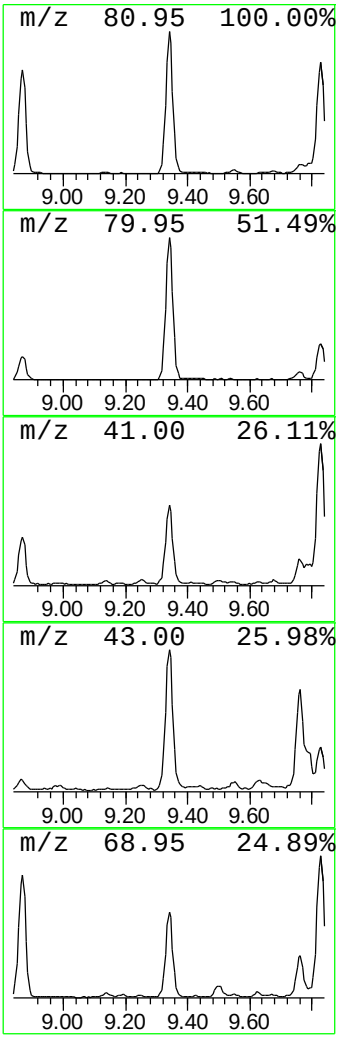
****

Peak Number 3 Bicyclo[2.2.1]heptan-2-ol, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.34	8.06 ng/ul	968366	Naphthalene-d8	10.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-ol, 1,3,3...	154	C10H18O	001632-73-1	96
2	Fenchol, exo-	154	C10H18O	022627-95-8	91
3	Bicyclo[2.2.1]heptan-2-ol, 1,3,3...	154	C10H18O	001632-73-1	90
4	Bicyclo[2.2.1]heptan-2-ol, 1,3,3...	154	C10H18O	001632-73-1	90
5	Bicyclo[2.2.1]heptan-2-ol, 1,3,3...	154	C10H18O	001632-73-1	87





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

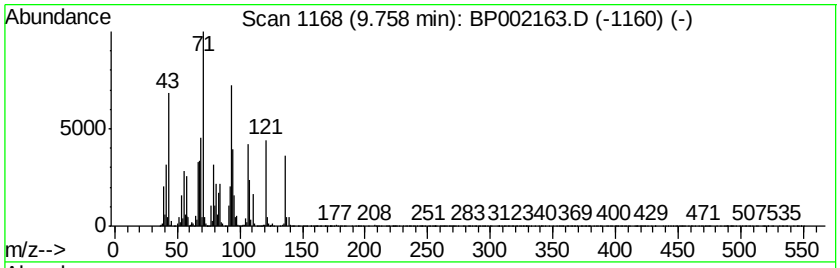
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

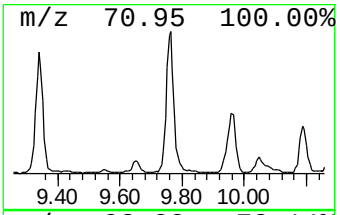
****

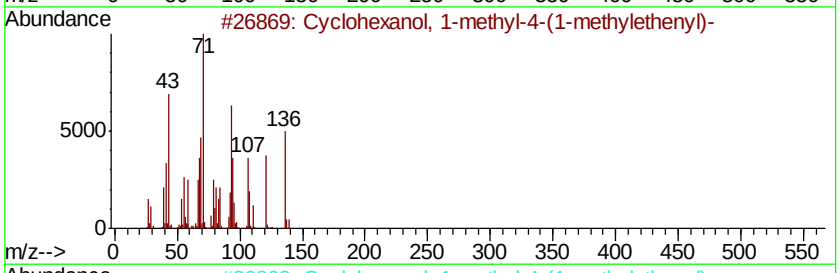
Peak Number 4 Cyclohexanol, 1-methyl-4-(1... Concentration Rank 7

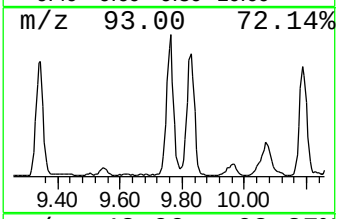
R.T.	EstConc	Area	Relative to ISTD	R.T.
9.76	4.07 ng/ul	488551	Naphthalene-d8	10.40

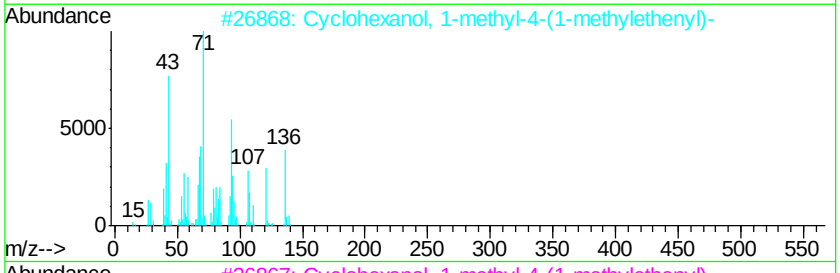
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanol, 1-methyl-4-(1-meth...	154	C10H18O	000138-87-4	94
2	Cyclohexanol, 1-methyl-4-(1-meth...	154	C10H18O	000138-87-4	91
3	Cyclohexanol, 1-methyl-4-(1-meth...	154	C10H18O	000138-87-4	90
4	Cyclohexanol, 1-methyl-4-(1-meth...	154	C10H18O	000138-87-4	64
5	p-Menth-8-en-1-ol, stereoisomer	154	C10H18O	007299-40-3	58

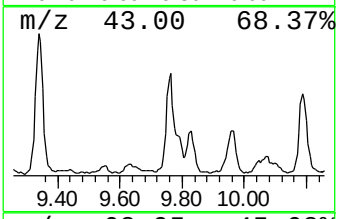


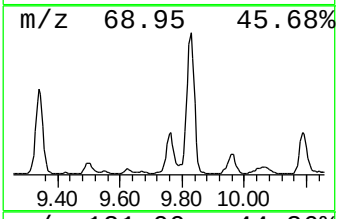


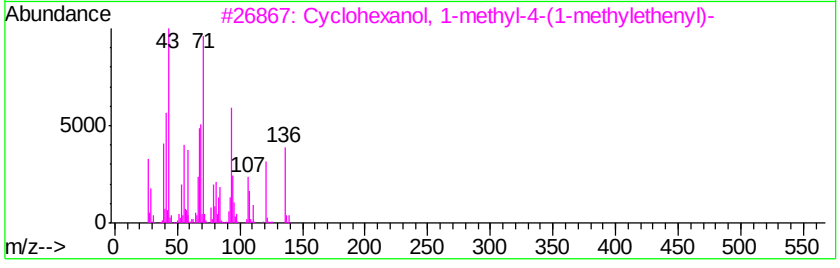


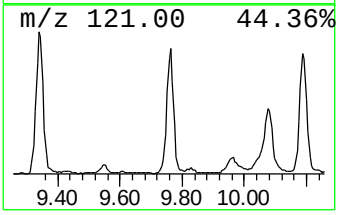












Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

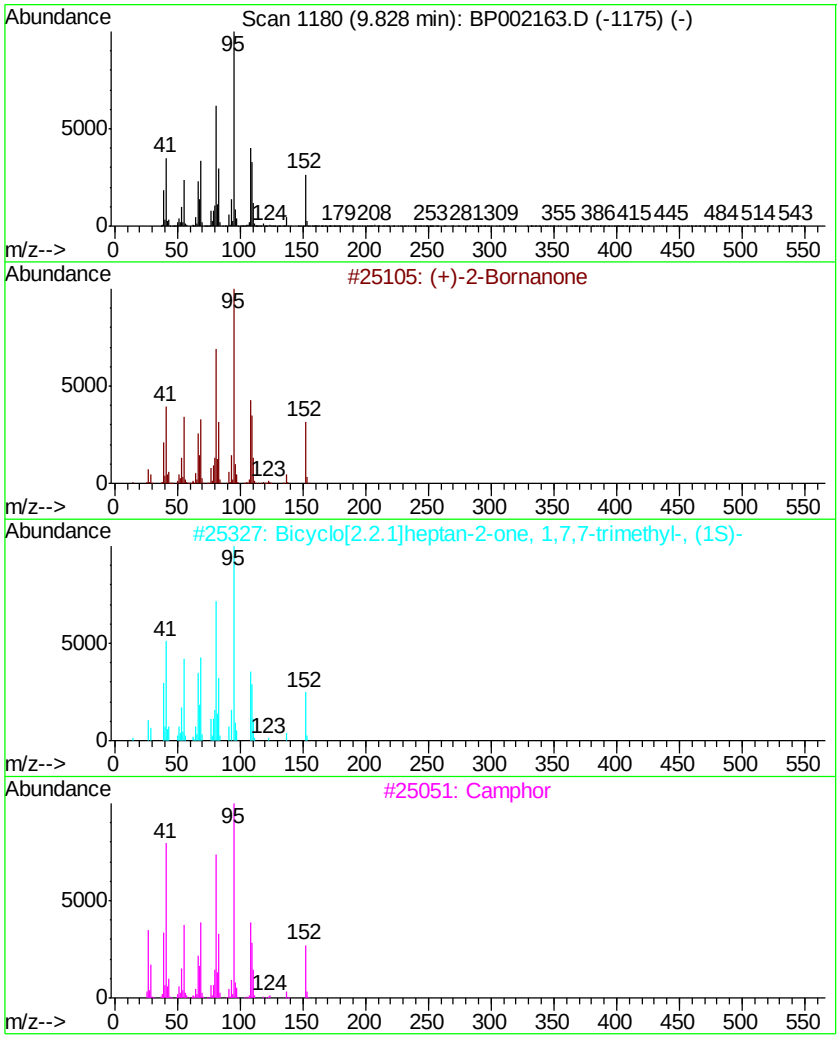
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

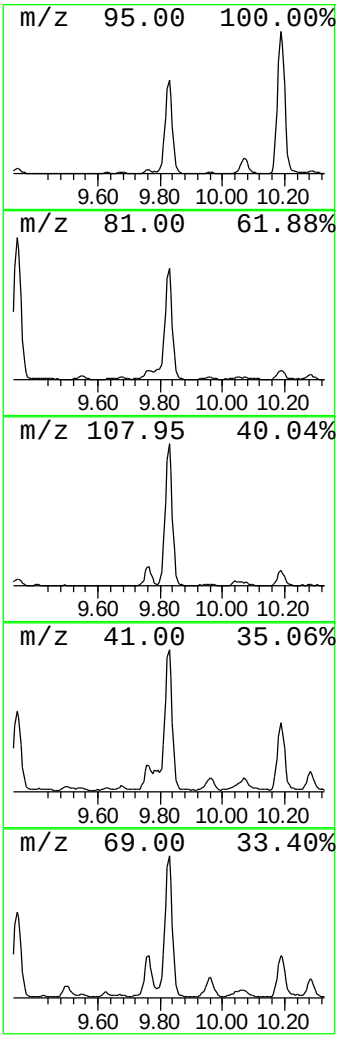
****

Peak Number 5 (+)-2-Bornanone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.83	10.26 ng/ul	1231670	Naphthalene-d8	10.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-2-Bornanone	152	C10H16O	000464-49-3	98
2	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	97
3	Camphor	152	C10H16O	000076-22-2	96
4	(+)-2-Bornanone	152	C10H16O	000464-49-3	96
5	(+)-2-Bornanone	152	C10H16O	000464-49-3	96





Library Search Compound Report

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Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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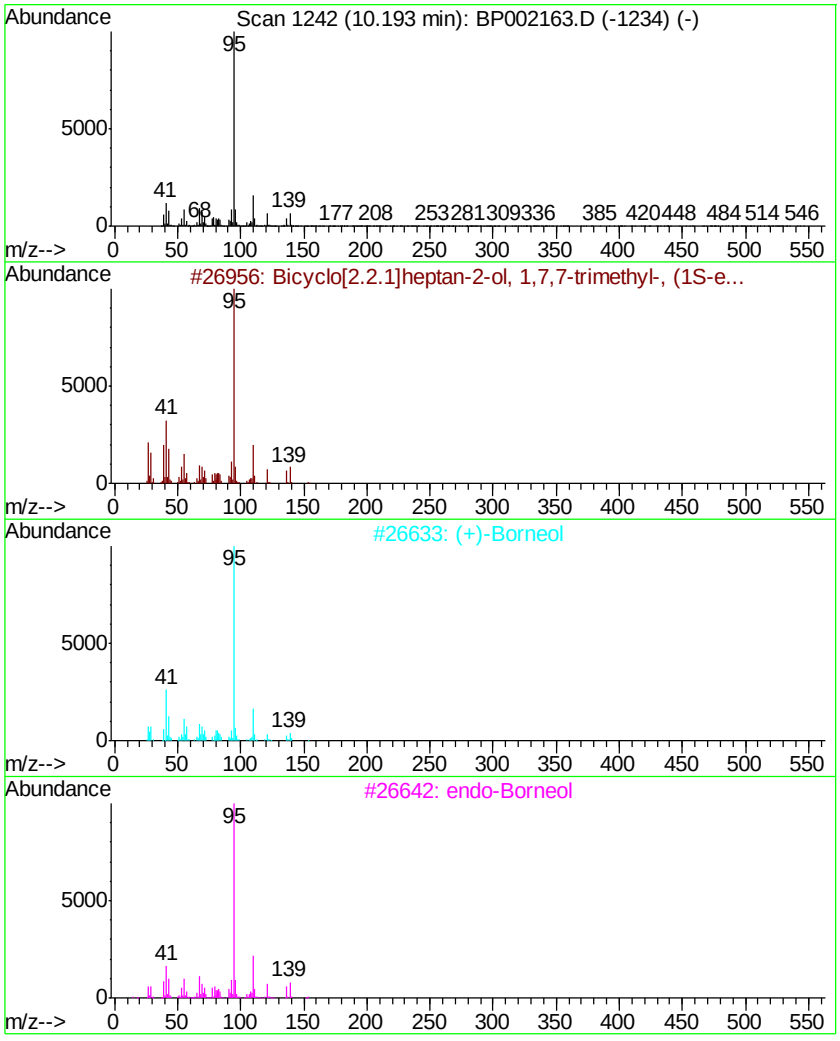
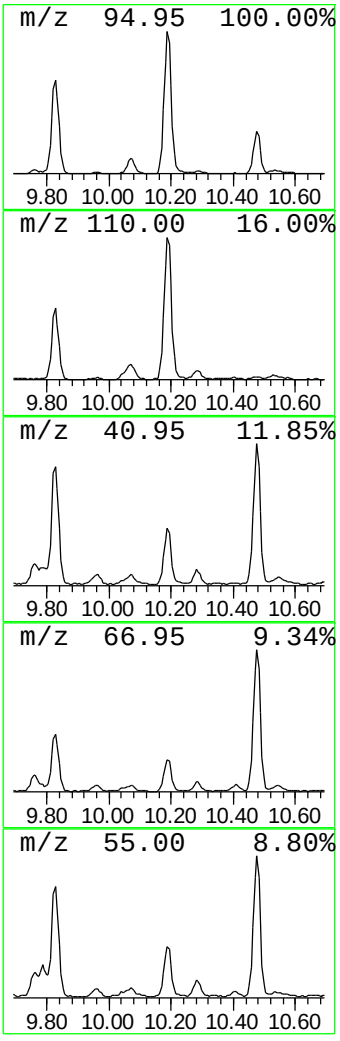
Peak Number 6 Bicyclo[2.2.1]heptan-2-ol, ... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.19	7.14 ng/ul	856968	Naphthalene-d8	10.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-ol, 1,7,7...	154	C10H18O	000464-45-9	91
2	(+)-Borneol	154	C10H18O	1000150-76-3	91
3	endo-Borneol	154	C10H18O	000507-70-0	91
4	2-Amino-4-methylbut-2-enenitrile	110	C6H10N2	1000197-39-9	64
5	endo-Borneol	154	C10H18O	000507-70-0	64





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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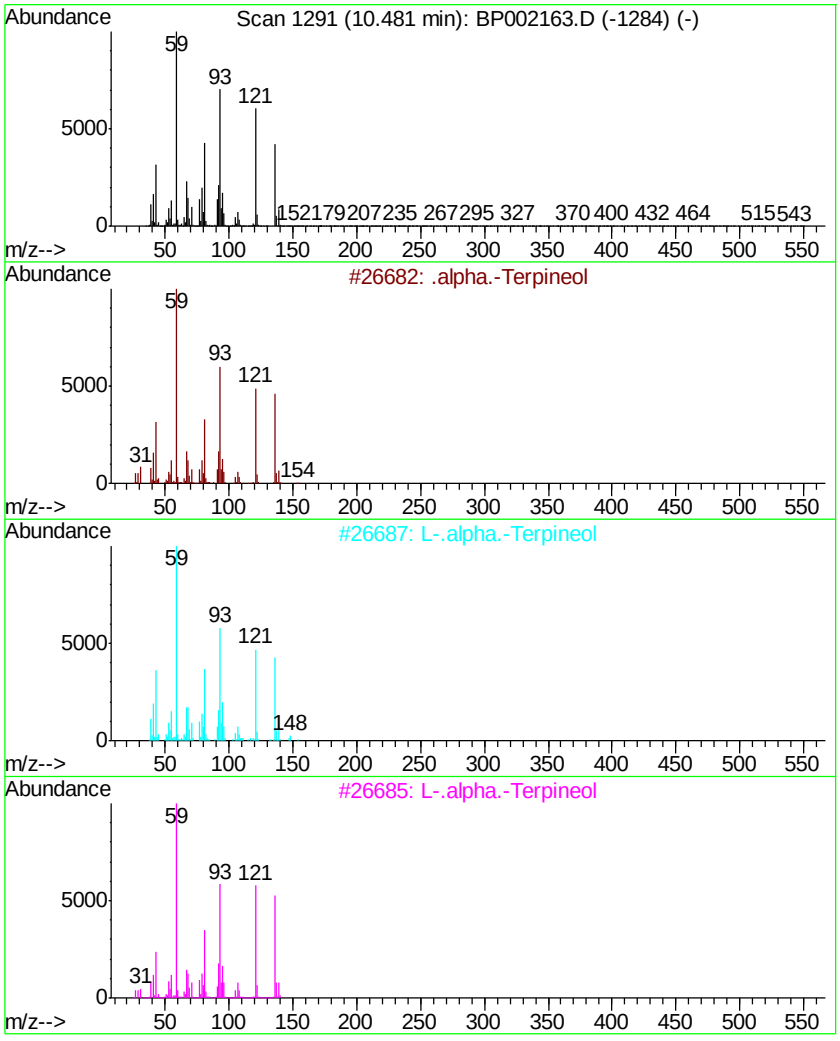
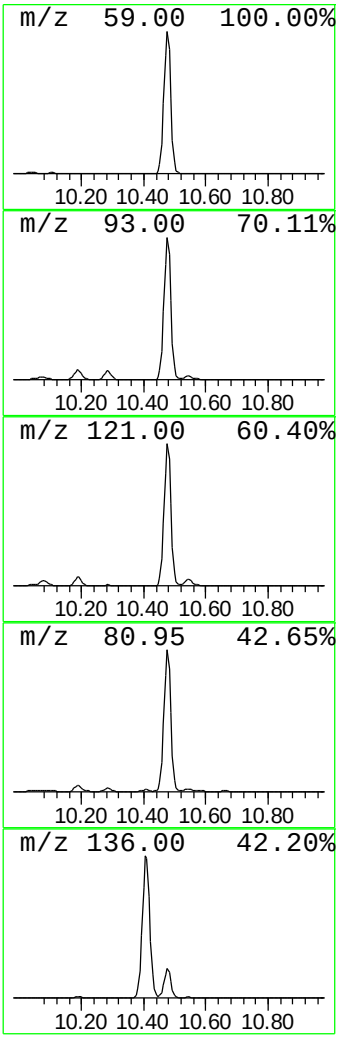
Peak Number 7 .alpha.-Terpineol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.48	27.19 ng/ul	3265250	Naphthalene-d8	10.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Terpineol	154	C10H18O	000098-55-5	83
2	L-.alpha.-Terpineol	154	C10H18O	010482-56-1	72
3	L-.alpha.-Terpineol	154	C10H18O	010482-56-1	72
4	1,3-Cyclohexadiene, 1-methyl-4-(...	136	C10H16	000099-86-5	50
5	(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	50





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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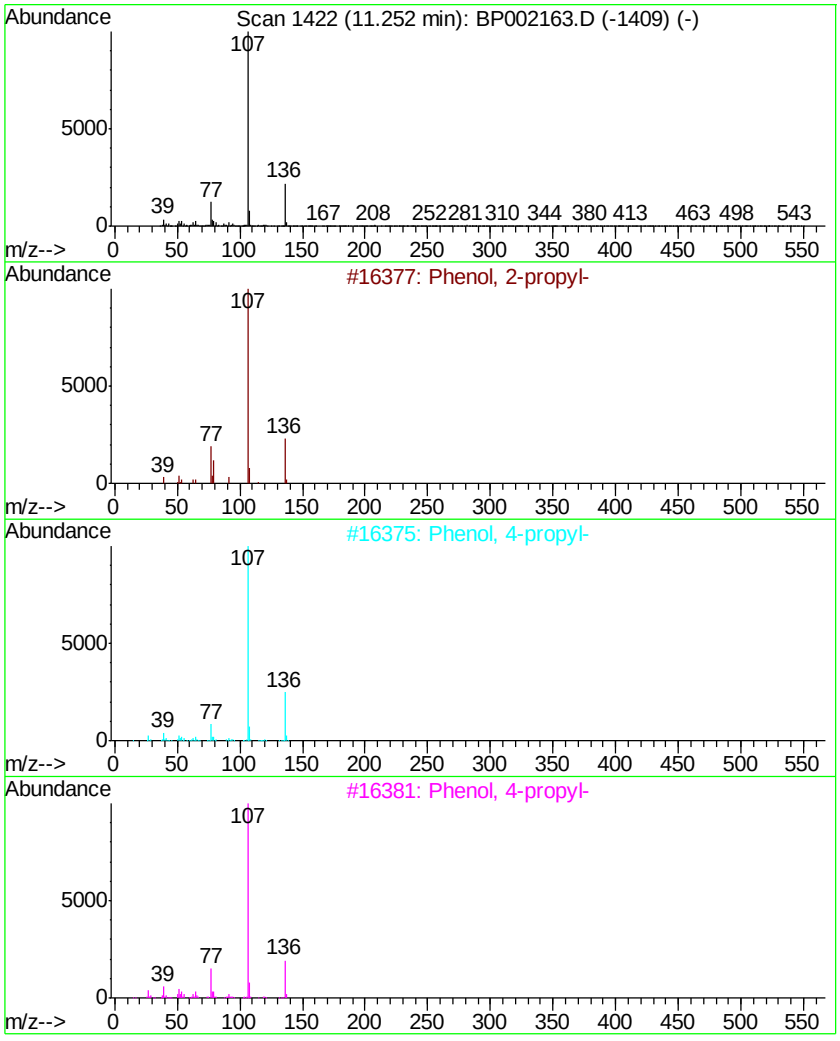
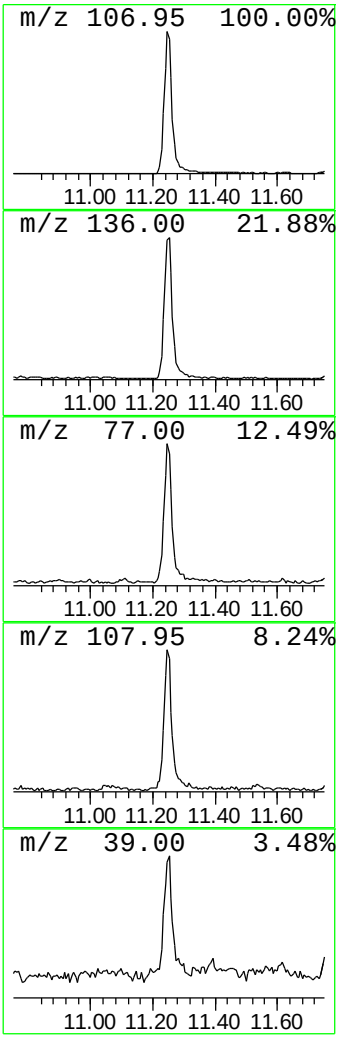
Peak Number 8 Phenol, 2-propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.25	3.02 ng/ul	363008	Naphthalene-d8	10.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-propyl-	136	C9H12O	000644-35-9	90
2	Phenol, 4-propyl-	136	C9H12O	000645-56-7	90
3	Phenol, 4-propyl-	136	C9H12O	000645-56-7	87
4	Formic acid, 2-propylphenyl ester	164	C10H12O2	1000368-96-6	64
5	Phenol, 2-propyl-	136	C9H12O	000644-35-9	60





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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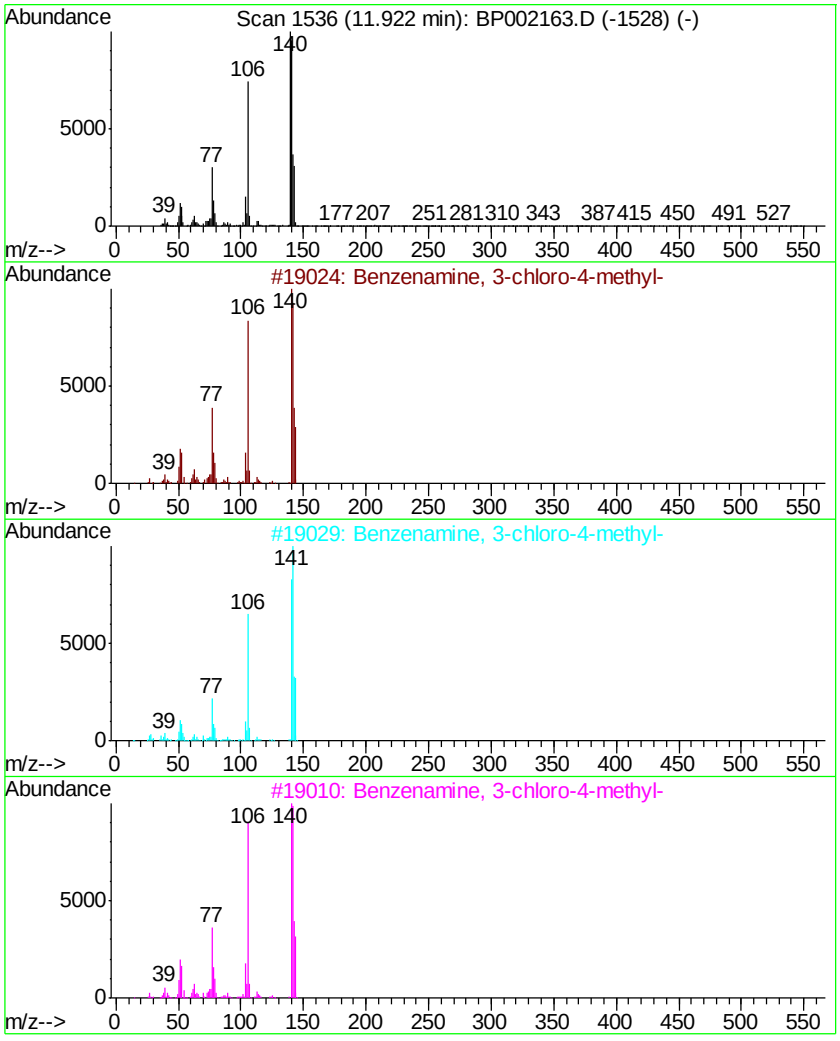
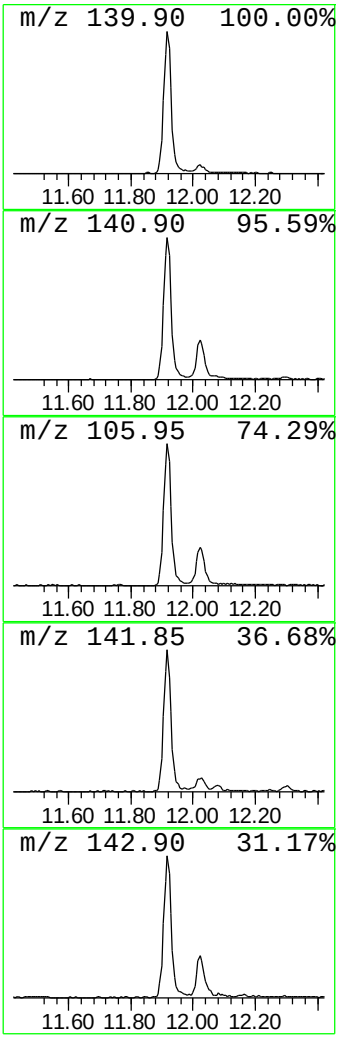
Peak Number 9 Benzenamine, 3-chloro-4-met... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	3.39 ng/ul	406883	Naphthalene-d8	10.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	98
2	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	96
3	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	96
4	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	95
5	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	94





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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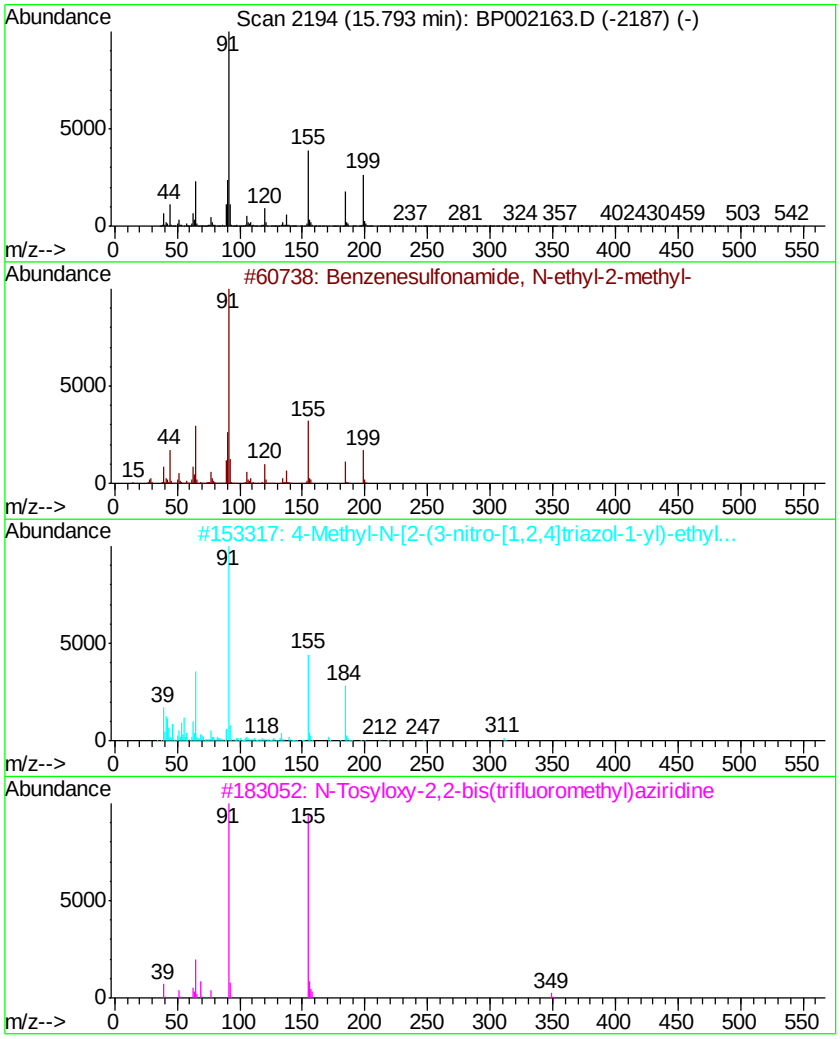
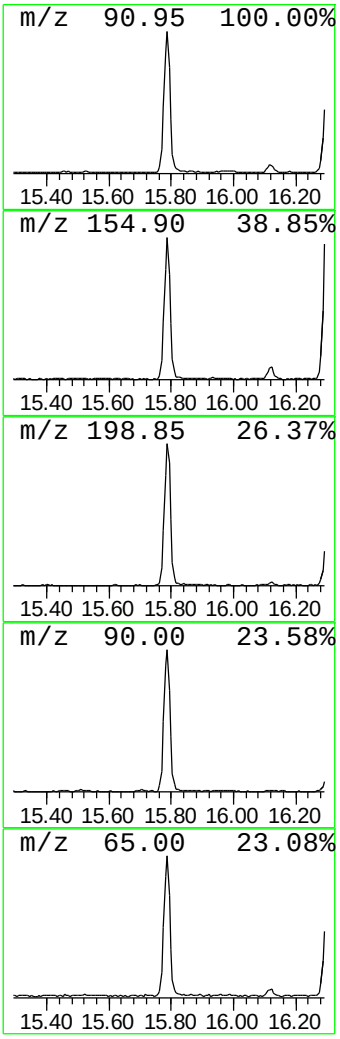
Peak Number 10 Benzenesulfonamide, N-ethyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.79	2.03 ng/ul	437417	Phenanthrene-d10	17.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	96
2	4-Methyl-N-[2-(3-nitro-[1,2,4]tr...	311	C11H13N5O4S	1000301-03-5	50
3	N-Tosyloxy-2,2-bis(trifluorometh...	349	C11H9F6NO3S	038436-38-3	49
4	Benzene, 1-[(chloromethyl)sulfon...	204	C8H9ClO2S	007569-26-8	47
5	Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	47





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP031020\
Data File : BP002163.D
Acq On : 10 Mar 2020 22:29
Operator : CG/JU
Sample : L1843-08DL 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB5T3DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzenamine, 3-me...	8.61	20.1	ng/ul	1672400	1	7.63	1662730	20.0
Bicyclo[2.2.1]hep...	8.87	4.1	ng/ul	339697	1	7.63	1662730	20.0
Bicyclo[2.2.1]hep...	9.34	8.1	ng/ul	968366	2	10.40	2401810	20.0
Cyclohexanol, 1-m...	9.76	4.1	ng/ul	488551	2	10.40	2401810	20.0
(+)-2-Bornanone	9.83	10.3	ng/ul	1231670	2	10.40	2401810	20.0
Bicyclo[2.2.1]hep...	10.19	7.1	ng/ul	856968	2	10.40	2401810	20.0
.alpha.-Terpineol	10.48	27.2	ng/ul	3265250	2	10.40	2401810	20.0
Phenol, 2-propyl-	11.25	3.0	ng/ul	363008	2	10.40	2401810	20.0
Benzenamine, 3-ch...	11.92	3.4	ng/ul	406883	2	10.40	2401810	20.0
Benzenesulfonamid...	15.79	2.0	ng/ul	437417	4	17.02	4315610	20.0