LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004372.D
Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

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-BP121820MA.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	3.352	24	28	36	rVB	303847	423319	10.86%	0.825%
2	4.116	154	158	162	rBV	65867	85628	2.20%	0.167%
3	4.481	216	220	232	rVB	115203	191246	4.91%	0.373%
4	6.104	493	496	502	rVB	17367	28868	0.74%	0.056%
5	6.992	639	647	649	rBV2	138994	257356	6.60%	0.501%
6	7.157	669	675	685	rBV	325360	526784	13.51%	1.026%
7	7.357	702	709	716	rBV	405094	649839	16.67%	1.266%
8	7.416	716	719	727	rVB	68695	108029	2.77%	0.211%
9	7.822	780	788	795	rBV	778968	1282462	32.90%	2.499%
10	7.887	795	799	805	rVB3	29753	49843	1.28%	0.097%
11	8.204	847	853	860	rVB4	18155	39756	1.02%	0.077%
12	8.528	901	908	916	rVV	361843	581682	14.92%	1.133%
13	8.722	935	941	948	rBV4	22356	47971	1.23%	0.093%
14	8.804	948	955	965	rBV	1947276	3168439	81.27%	6.174%
15	8.922	970	975	979	rVV3	35088	58148	1.49%	0.113%
16	8.981	979	985	995	rVV	380532	677953	17.39%	1.321%
17	9.063	995	999	1003	rVB2	22984	34181	0.88%	0.067%
18	9.181	1009	1019	1026	rVB9	20869	53030	1.36%	0.103%
19	9.339	1039	1046	1050	rBV5	34361	62596	1.61%	0.122%
20	9.428	1051	1061	1071	rVB2	124748	281477	7.22%	0.548%
21	9.622	1090	1094	1101	rVB8	12022	27178	0.70%	0.053%
22	9.698	1101	1107	1114	rBV	346680	591359	15.17%	1.152%
23	9.834	1120	1130	1134	rBV7	29268	64292	1.65%	0.125%
24	9.881	1134	1138	1144	rVV	32521	57172	1.47%	0.111%
25	9.928	1144	1146	1154	rVB8	10557	20401	0.52%	0.040%
26	10.034	1154	1164	1171	rVB2	78282	146033	3.75%	0.285%
27	10.110	1171	1177	1179	rBV4	11292	20896	0.54%	0.041%
28	10.163	1179	1186	1192	rVB2	75545	158958	4.08%	0.310%
29	10.239	1192	1199	1218	rVB	621037	1183496	30.36%	2.306%
30	10.381	1218	1223	1224	rBV5	16485	20385	0.52%	0.040%
31	10.533	1237	1249	1257	rBV2	41377	84505	2.17%	0.165%
32	10.622	1257	1264	1270	rVV	1089687	1897984	48.68%	3.698%
33	10.669	1270	1272	1277	rVV2	25720	35448	0.91%	0.069%
34	10.751	1277	1286	1299	rVB	557403	1050112	26.94%	2.046%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
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Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

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-BP121820MA.M
Title : SVOA CALIBRATION

35	10.875	1299	1307	1316	rBV8	30753	92198	2.36%	0.180%
36	11.045	1329	1336	1342	rBV7	20733	49049	1.26%	0.096%
37	11.110	1342	1347	1354	rVB9	9498	22526	0.58%	0.044%
38	11.175	1354	1358	1361	rBV6	17508	31506	0.81%	0.061%
39	11.228	1365	1367	1372	rVB5	20191	25067	0.64%	0.049%
40	11.439	1395	1403	1404	rVV5	31797	67773	1.74%	0.132%
41	11.463	1404	1407	1413	rVV3	46530	89404	2.29%	0.174%
42	11.533	1413	1419	1423	rVV8	22262	56937	1.46%	0.111%
43	11.586	1423	1428	1434	rVV7	42247	88564	2.27%	0.173%
44	11.663	1434	1441	1450	rVB7	23737	70800	1.82%	0.138%
45	11.822	1464	1468	1474	rVB	30147	52937	1.36%	0.103%
46	11.963	1487	1492	1504	rVB	60447	127939	3.28%	0.249%
47	12.069	1504	1510	1513	rBV7	13887	26461	0.68%	0.052%
48	12.122	1513	1519	1527	rBV	758192	1246119	31.96%	2.428%
49	12.228	1532	1537	1544	rVB	146708	241492	6.19%	0.471%
50	12.339	1551	1556	1561	rBV7	20552	43243	1.11%	0.084%
51	12.504	1579	1584	1589	rBV8	17594	31026	0.80%	0.060%
52	12.575	1589	1596	1598	rBV8	12288	27351	0.70%	0.053%
53	12.622	1599	1604	1609	rVV2	25478	44839	1.15%	0.087%
54	12.722	1617	1621	1625	rVB5	16291	25447	0.65%	0.050%
55	12.845	1633	1642	1648	rBV10	19440	48131	1.23%	0.094%
56	13.180	1694	1699	1706	rVV8	22819	56166	1.44%	0.109%
57	13.263	1708	1713	1716	rVV5	20920	46223	1.19%	0.090%
58	13.310	1717	1721	1728	rVB6	34255	77221	1.98%	0.150%
59	13.369	1728	1731	1736	rBV7	18314	36697	0.94%	0.072%
60	13.504	1750	1754	1760	rBV3	32361	72364	1.86%	0.141%
61	13.557	1760	1763	1768	rVB	36665	54283	1.39%	0.106%
62	13.698	1782	1787	1792	rBV9	17227	40329	1.03%	0.079%
63	13.874	1812	1817	1822	rBV	1054882	1484133	38.07%	2.892%
64	13.922	1822	1825	1836	rVB	552832	749605	19.23%	1.461%
65	14.022	1836	1842	1848	rBV8	35695	77466	1.99%	0.151%
66	14.163	1856	1866	1878	rBV	1190065	1847528	47.39%	3.600%
67	14.274	1880	1885	1890	rBV	86702	138761	3.56%	0.270%
68	14.392	1899	1905	1910	rBV4	109258	174242	4.47%	0.340%
69	14.469	1912	1918	1926	rVV2	1698500	2614517	67.06%	5.095%
70	14.539	1926	1930	1935	rVB4	20885	42611	1.09%	0.083%
71	14.674	1948	1953	1959	rBV	117269	185078	4.75%	0.361%

LSC Area Percent Report

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Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

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-BP121820MA.M
Title : SVOA CALIBRATION

72	15.216	2041	2045	2050	rBV	41693	84644	2.17%	0.165%
73	15.333	2062	2065	2069	rBV4	17858	23160	0.59%	0.045%
74	15.469	2082	2088	2094	rBV	1492795	2150199	55.15%	4.190%
75	15.592	2103	2109	2114	rBV	353812	523590	13.43%	1.020%
76	15.721	2126	2131	2143	rVB7	56986	165979	4.26%	0.323%
77	15.927	2163	2166	2169	rBV4	18150	21475	0.55%	0.042%
78	15.980	2170	2175	2185	rVB	492385	708892	18.18%	1.381%
79	16.151	2197	2204	2209	rBV2	183920	306931	7.87%	0.598%
80	16.204	2210	2213	2216	rVB4	17852	21901	0.56%	0.043%
81	16.316	2228	2232	2236	rBV2	27873	37034	0.95%	0.072%
82	16.368	2238	2241	2247	rVB5	18950	31931	0.82%	0.062%
83	16.498	2255	2263	2267	rBV	315091	487069	12.49%	0.949%
84	16.545	2267	2271	2277	rVB3	43925	81443	2.09%	0.159%
85	16.763	2305	2308	2314	rVB	59012	78294	2.01%	0.153%
86	17.233	2381	2388	2393	rBV	2132497	3152442	80.86%	6.143%
87	17.327	2398	2404	2415	rVB	1679093	2490325	63.88%	4.853%
88	18.127	2535	2540	2548	rBV	166029	301891	7.74%	0.588%
89	18.410	2585	2588	2592	rBV2	50304	65028	1.67%	0.127%
90	18.874	2663	2667	2676	rVB	117210	170230	4.37%	0.332%
91	19.362	2747	2750	2754	rBV6	32753	47715	1.22%	0.093%
92	19.568	2780	2785	2790	rBV	2229269	3103998	79.62%	6.048%
93	19.615	2790	2793	2799	rVB	170624	240621	6.17%	0.469%
94	21.039	3031	3035	3042	rBV2	668070	1084097	27.81%	2.112%
95	21.098	3042	3045	3052	rVB	343643	471680	12.10%	0.919%
96	21.321	3078	3083	3090	rBV	2865988	3696214	94.81%	7.202%
97	22.956	3359	3361	3367	rVB	144341	192634	4.94%	0.375%
98	23.527	3451	3458	3464	rBV	1560923	3129693	80.28%	6.099%
99	23.680	3477	3484	3497	rVB	1897249	3898528	100.00%	7.597%
100	24.274	3581	3585	3593	rVB2	192563	378427	9.71%	0.737%

Sum of corrected areas: 51318924

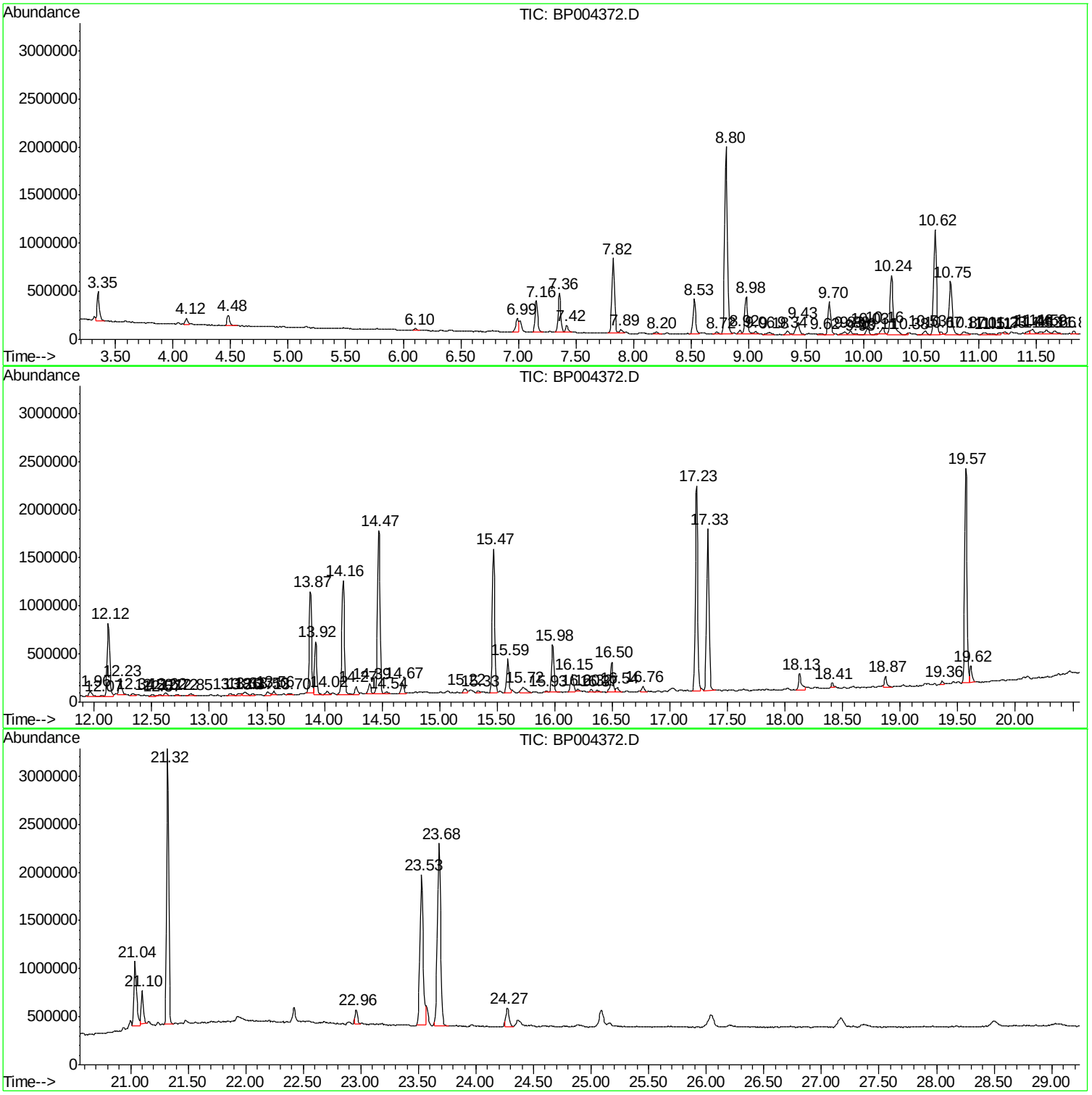
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004372.D
Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.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\BP121820\
Data File : BP004372.D
Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

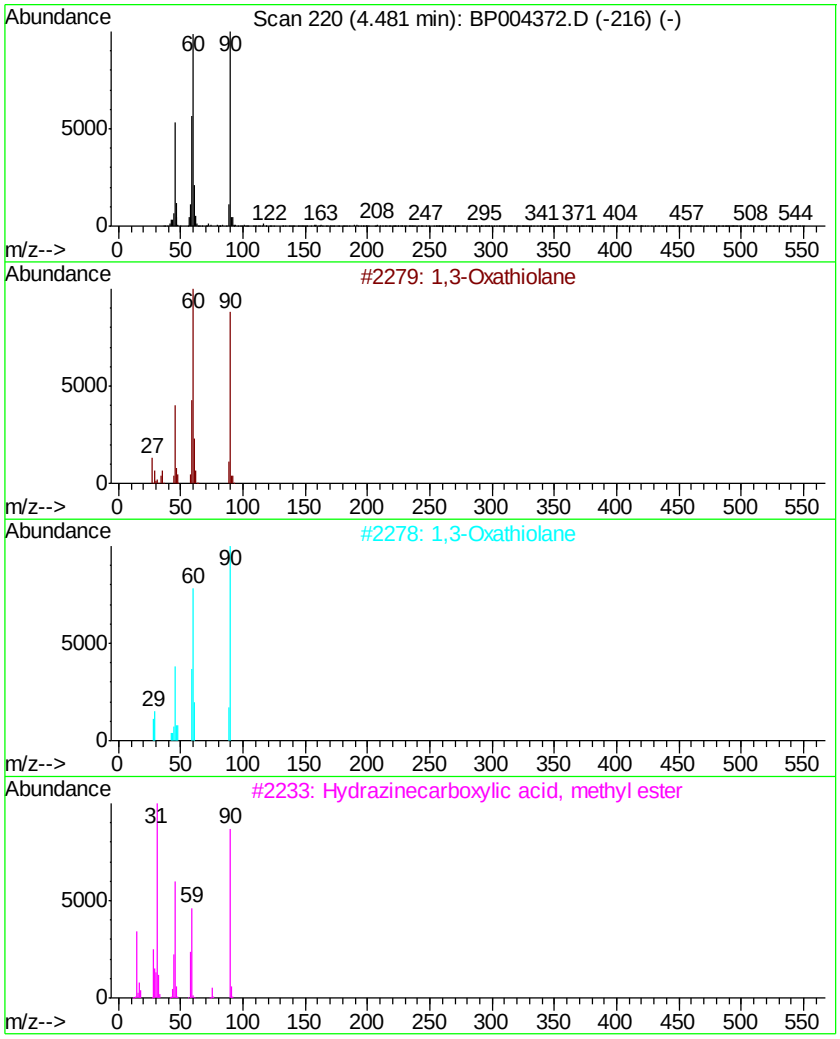
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

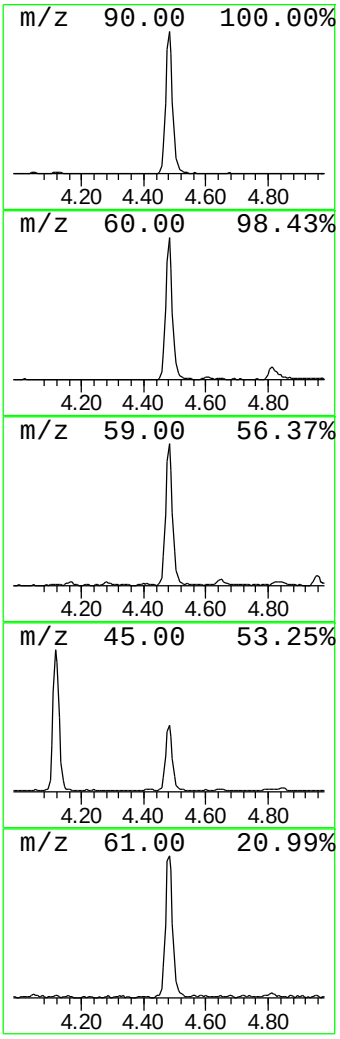
****

Peak Number 1 1,3-Oxathiolane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.48	2.98 ng/ul	191246	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Oxathiolane	90	C3H6OS	002094-97-5	90
2	1,3-Oxathiolane	90	C3H6OS	002094-97-5	78
3	Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	59
4	Heptanoic acid, 2-hydroxy-, meth...	160	C8H16O3	054340-91-9	53
5	Thiourea, methyl-	90	C2H6N2S	000598-52-7	50





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004372.D
Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

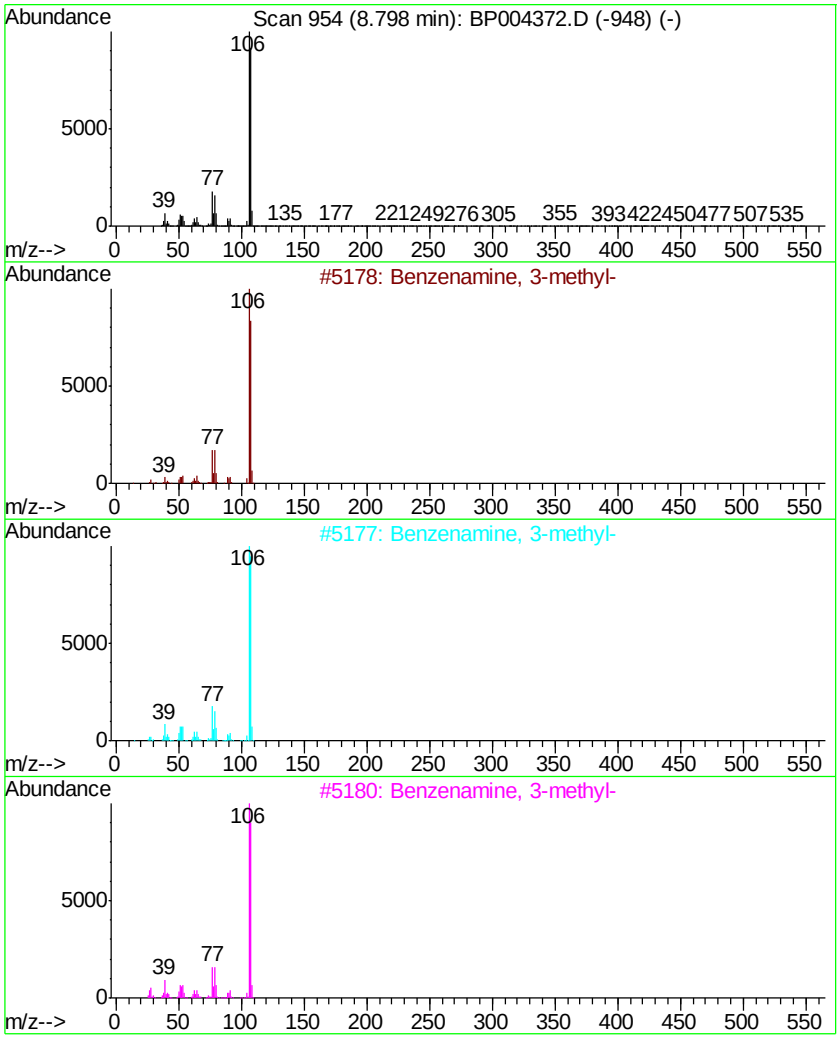
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

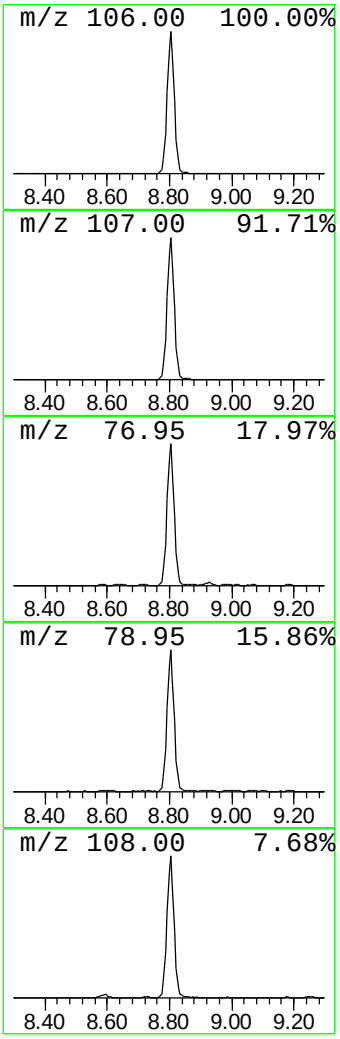
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.80	49.41 ng/ul	3168440	1,4-Dichlorobenzene-d4	7.82

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	o-Toluidine	107	C7H9N	000095-53-4	94





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004372.D
Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

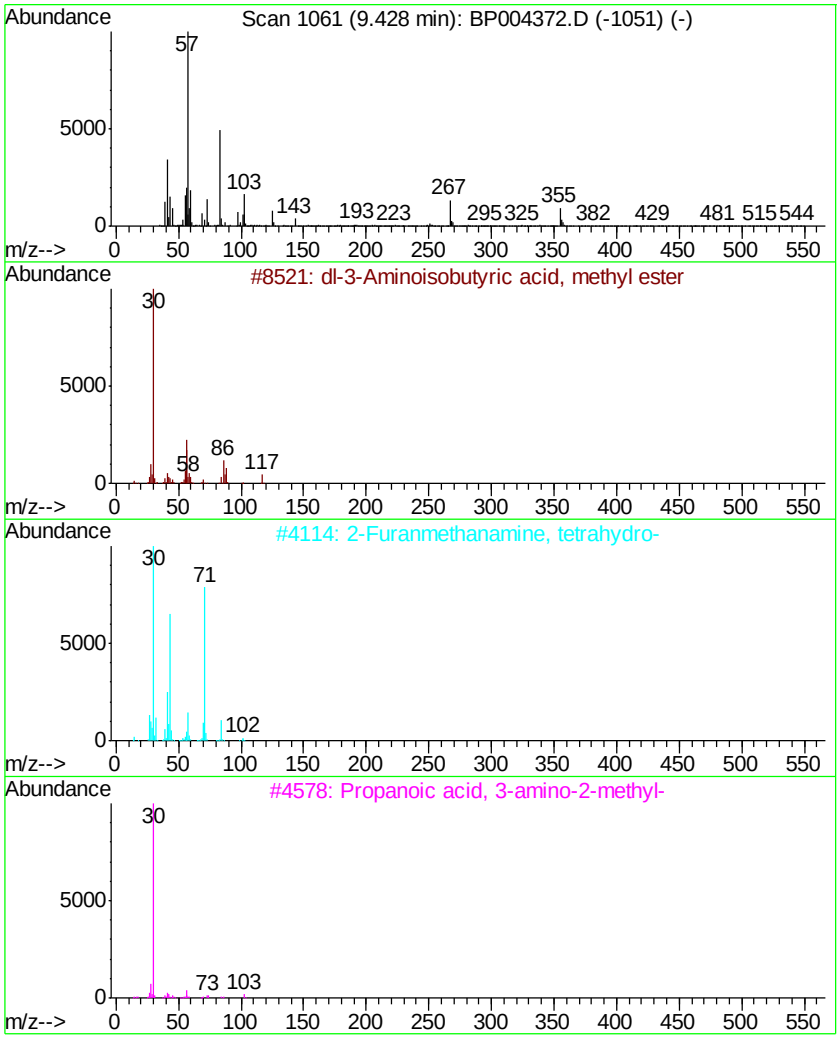
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

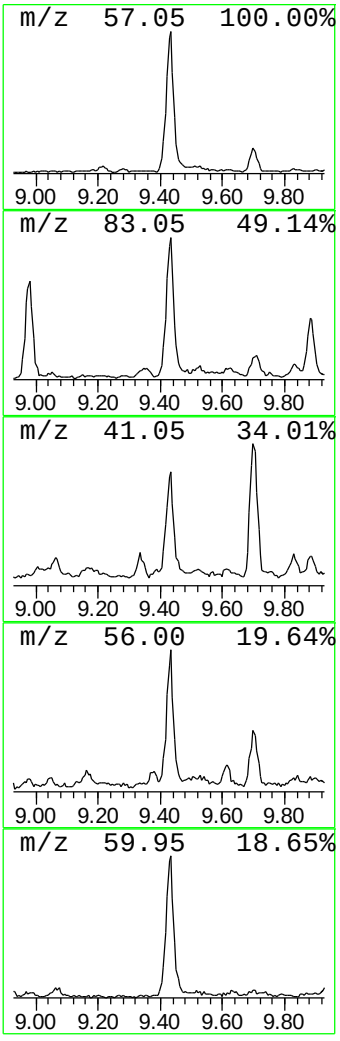
****

Peak Number 3 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.43	2.97 ng/ul	281477	Naphthalene-d8	10.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	dl-3-Aminoisobutyric acid, methy...	117	C5H11NO2	1000332-88-1	9
2	2-Furanmethanamine, tetrahydro-	101	C5H11NO	004795-29-3	7
3	Propanoic acid, 3-amino-2-methyl-	103	C4H9NO2	000144-90-1	4
4	1,3-Propanediol, 2-(hydroxymethy...	151	C4H9NO5	000126-11-4	4
5	L-Alanine, 3-[(aminocarbonyl)ami...	147	C4H9N3O3	001483-07-4	4





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004372.D
Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

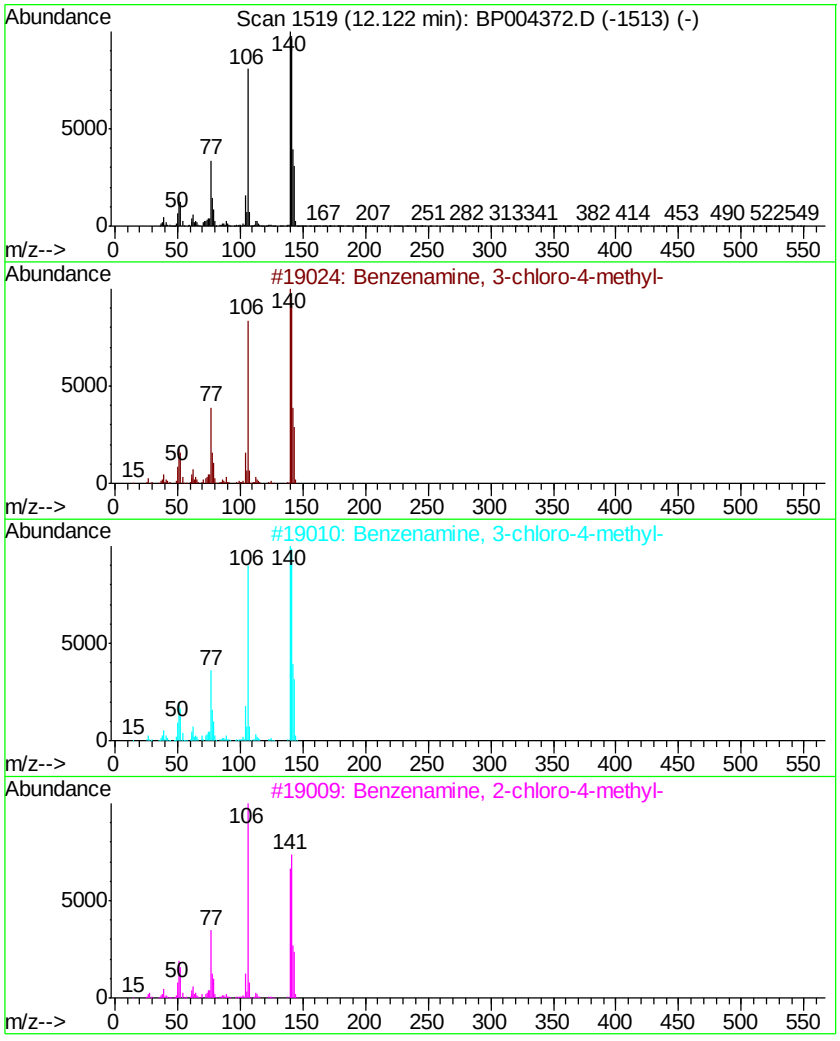
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

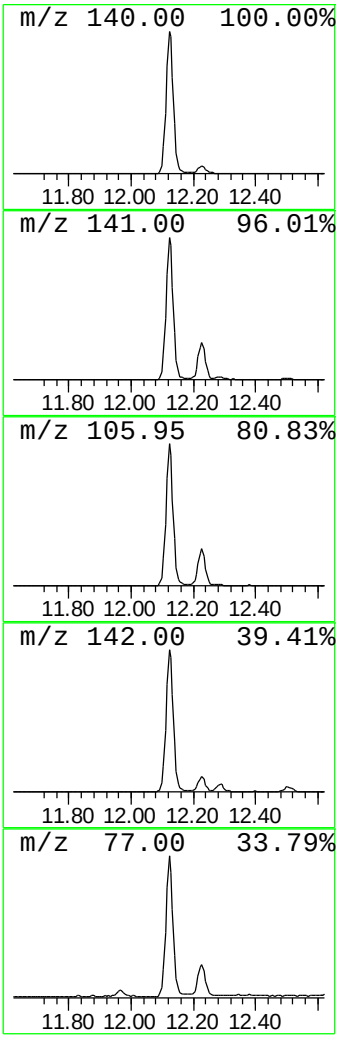
****

Peak Number 4 Benzenamine, 3-chloro-4-met... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.12	13.13 ng/ul	1246120	Naphthalene-d8	10.62

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	98
3	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	96
4	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	96
5	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	94





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004372.D
Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

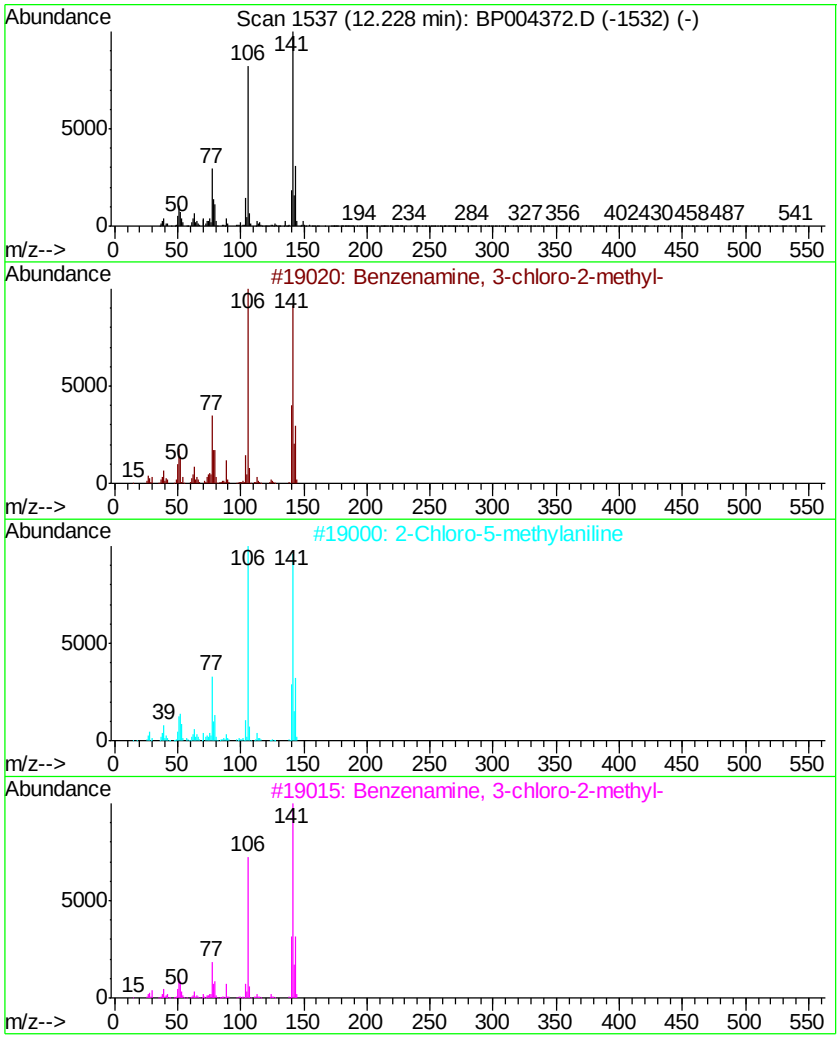
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

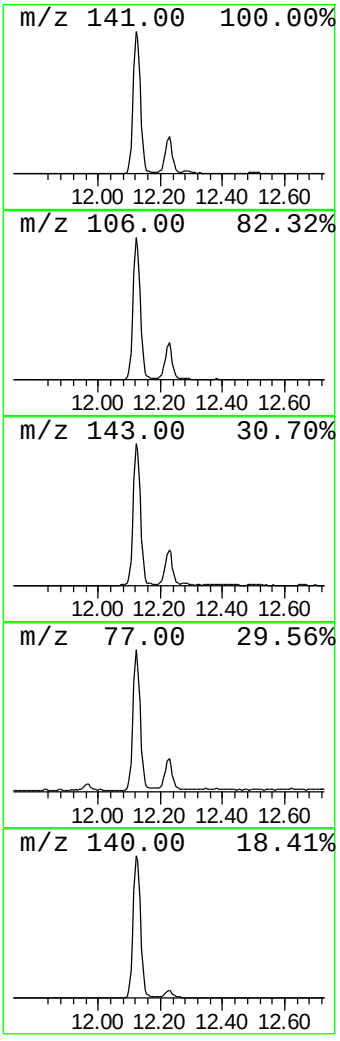
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzenamine, 3-chloro-2-met... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.23	2.54 ng/ul	241492	Naphthalene-d8	10.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-chloro-2-methyl-	141	C7H8ClN	000087-60-5	93
2	2-Chloro-5-methylaniline	141	C7H8ClN	000095-81-8	93
3	Benzenamine, 3-chloro-2-methyl-	141	C7H8ClN	000087-60-5	90
4	Benzenamine, 4-chloro-2-methyl-	141	C7H8ClN	000095-69-2	90
5	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	90





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Data File : BP004372.D
Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

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

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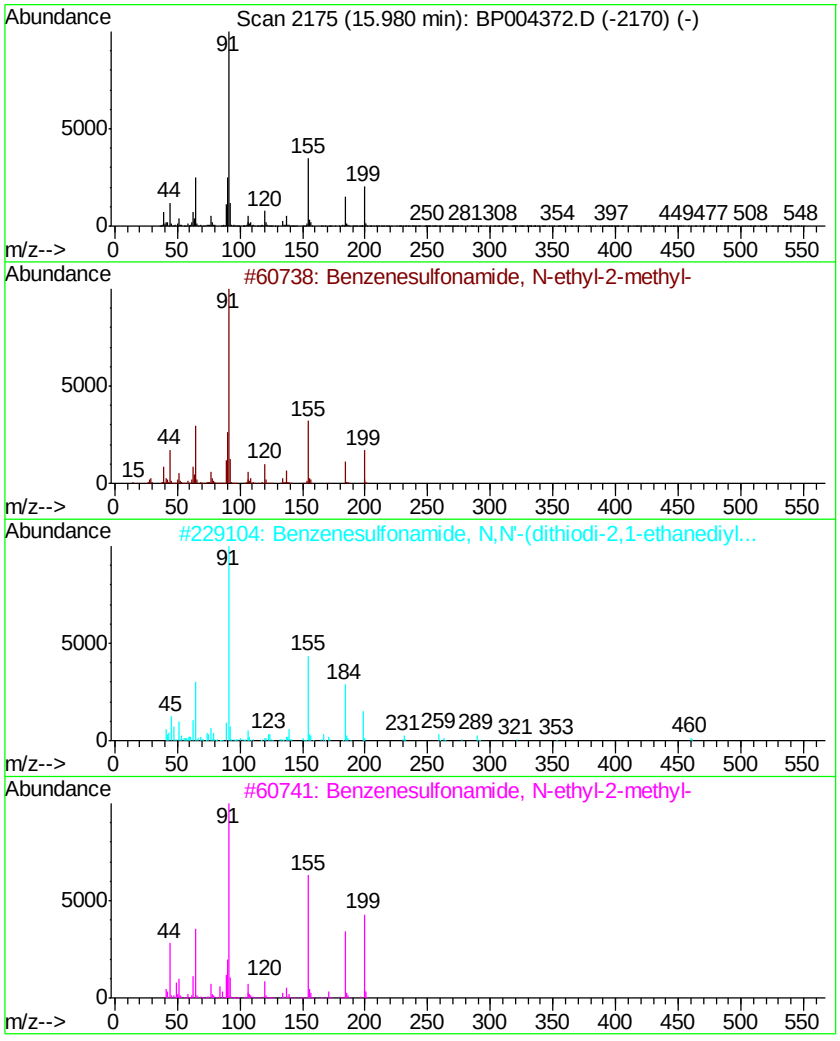
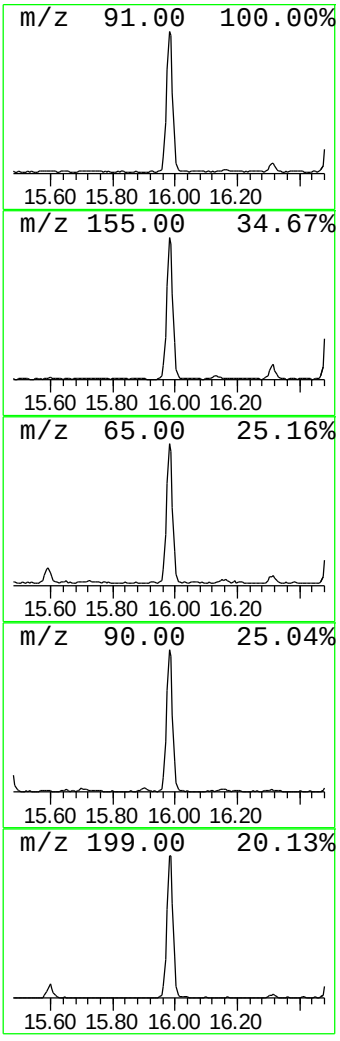
Peak Number 6 Benzenesulfonamide, N-ethyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.98	4.50 ng/ul	708892	Phenanthrene-d10	17.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	97
2	Benzenesulfonamide, N,N'-(dithio...	460	C18H24N2O4S4	023516-74-7	53
3	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	52
4	Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	49
5	Tolbutamide	270	C12H18N2O3S	000064-77-7	47





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Data File : BP004372.D
Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

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

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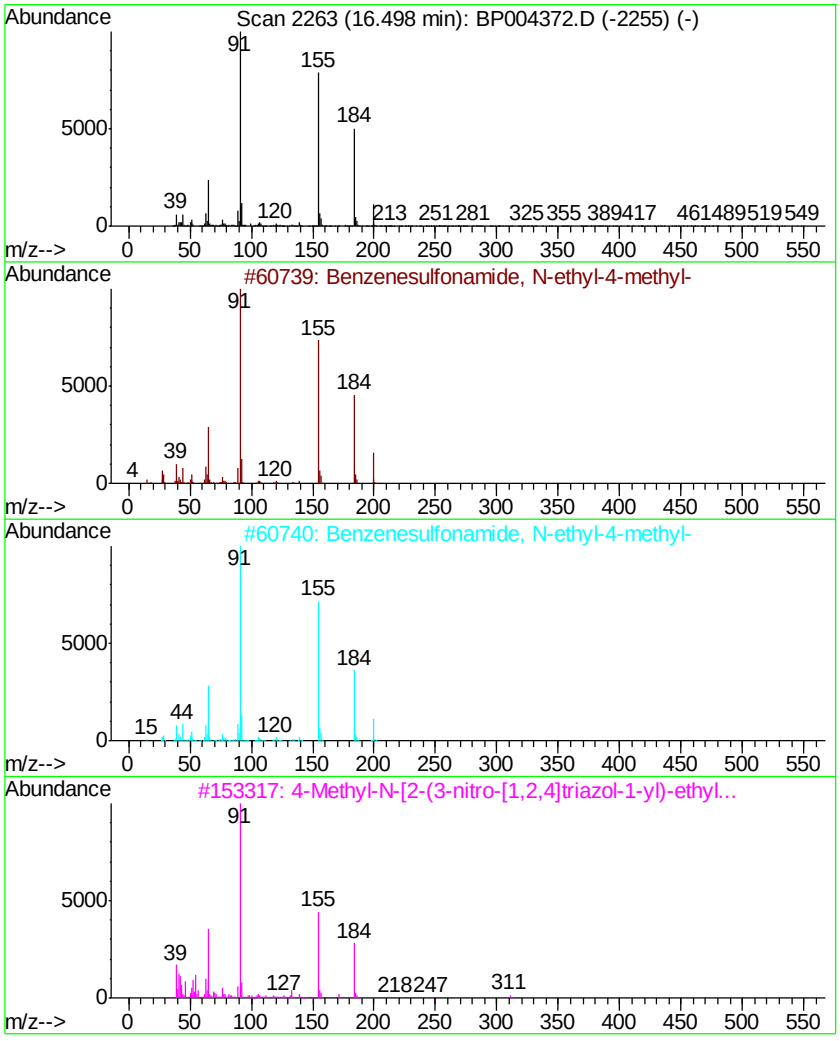
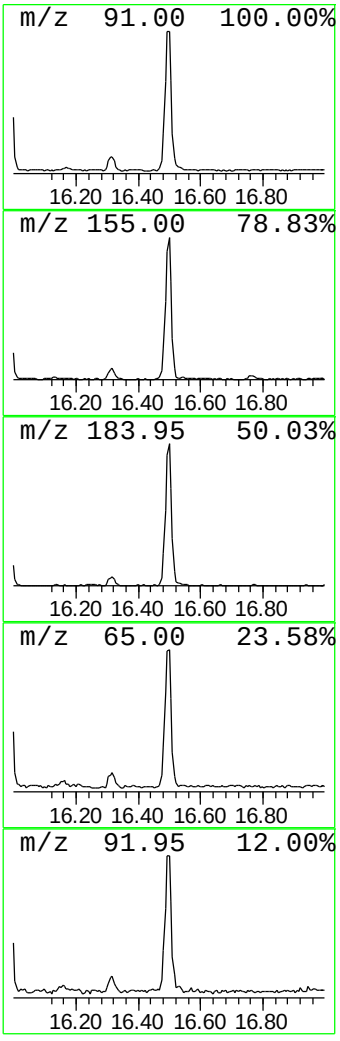
Peak Number 7 Benzenesulfonamide, N-ethyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	3.09 ng/ul	487069	Phenanthrene-d10	17.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	97
2	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	96
3	4-Methyl-N-[2-(3-nitro-[1,2,4]tr...	311	C11H13N5O4S	1000301-03-5	94
4	(Toluene-4-sulfonylamino)-acetic...	283	C13H17NO4S	1000190-48-0	72
5	Furazan-3-carboxylic acid, 4-ami...	325	C12H15N5O4S	1000304-69-4	64





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Data File : BP004372.D
Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

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

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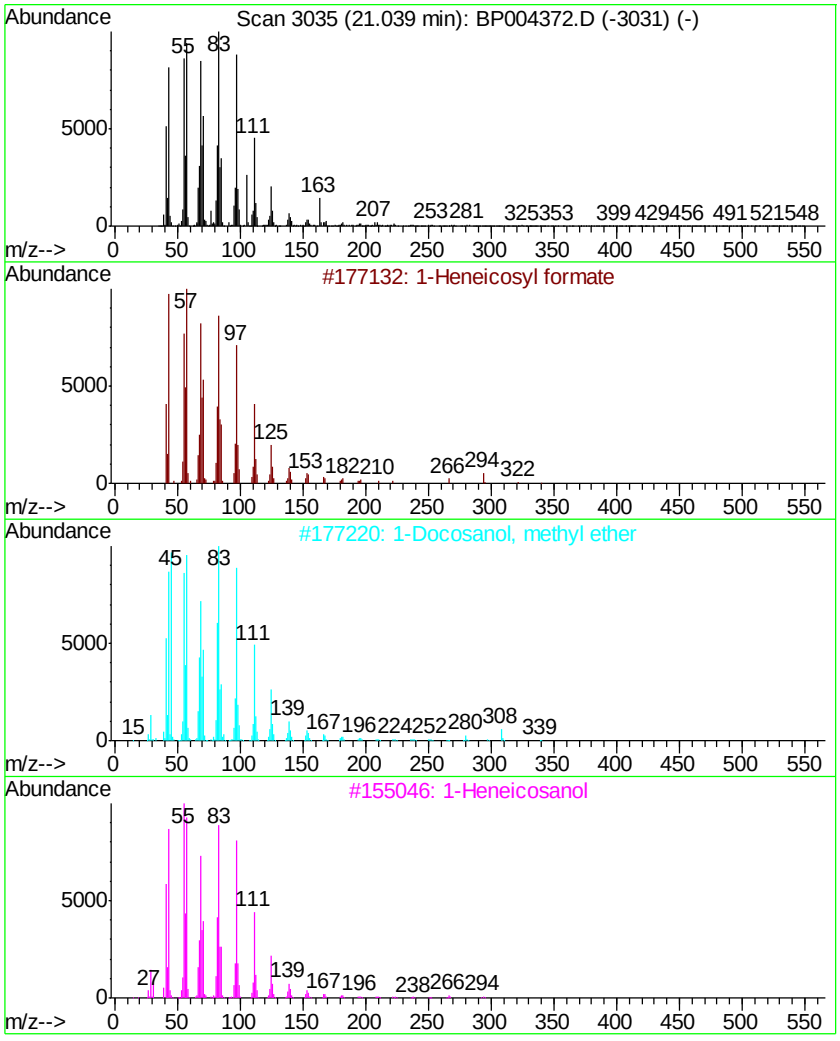
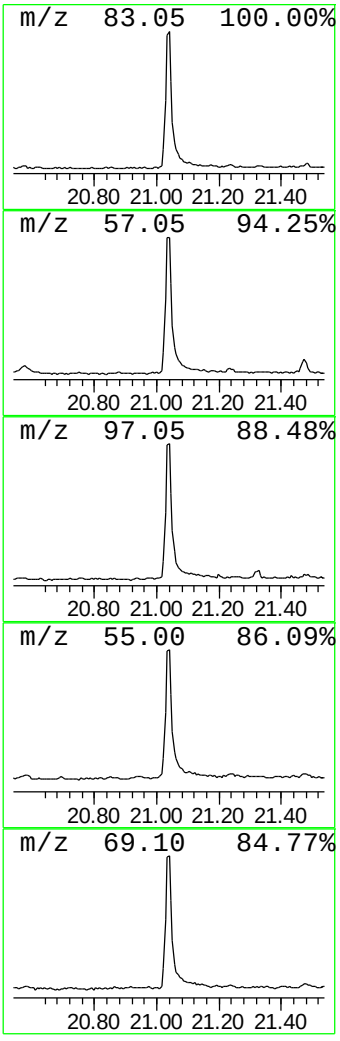
Peak Number 8 1-Heneicosyl formate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.04	5.87 ng/ul	1084100	Chrysene-d12	21.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	97
2	1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	96
3	1-Heneicosanol	312	C21H44O	015594-90-8	95
4	Cetene	224	C16H32	000629-73-2	94
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	94





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Data File : BP004372.D
Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

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

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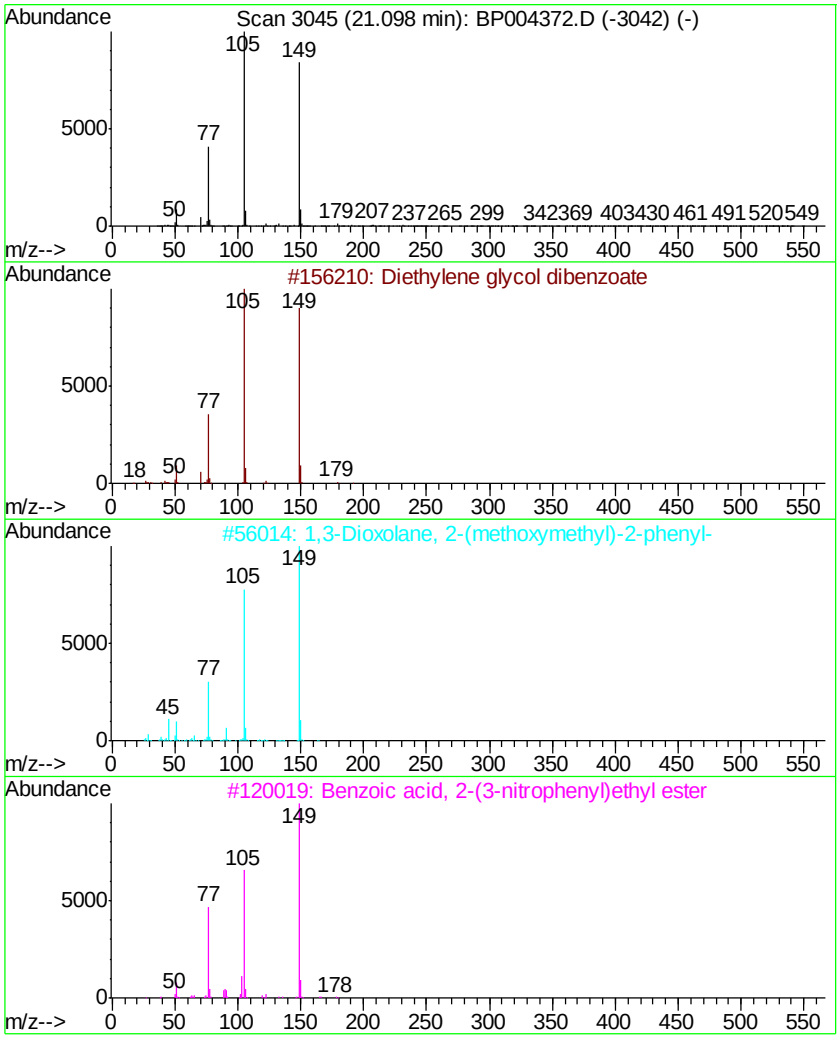
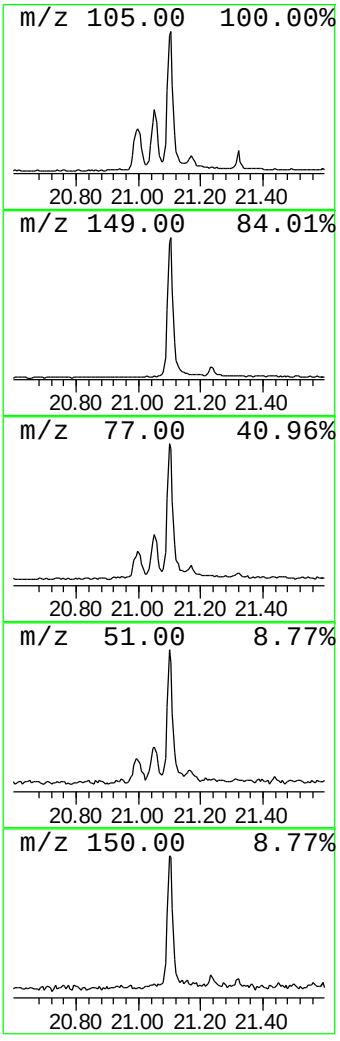
Peak Number 9 Diethylene glycol dibenzoate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.10	2.55 ng/ul	471680	Chrysene-d12	21.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	90
2	1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	78
3	Benzoic acid, 2-(3-nitrophenyl)e...	271	C15H13NO4	1000368-22-4	72
4	Benzoic acid, 2-(2-chlorophenoxy...	276	C15H13ClO3	1000368-25-9	64
5	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP121820\
Data File : BP004372.D
Acq On : 18 Dec 2020 23:40
Operator : CG/JU
Sample : L5128-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8E0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.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
1,3-Oxathiolane	4.48	3.0	ng/ul	191246	1	7.82	1282460	20.0
Benzenamine, 3-me...	8.80	49.4	ng/ul	3168440	1	7.82	1282460	20.0
unknown-01	9.43	3.0	ng/ul	281477	2	10.62	1897980	20.0
Benzenamine, 3-ch...	12.12	13.1	ng/ul	1246120	2	10.62	1897980	20.0
Benzenamine, 3-ch...	12.23	2.5	ng/ul	241492	2	10.62	1897980	20.0
Benzenesulfonamid...	15.98	4.5	ng/ul	708892	4	17.23	3152440	20.0
Benzenesulfonamid...	16.50	3.1	ng/ul	487069	4	17.23	3152440	20.0
1-Heneicosyl formate	21.04	5.9	ng/ul	1084100	5	21.32	3696210	20.0
Diethylene glycol...	21.10	2.5	ng/ul	471680	5	21.32	3696210	20.0