

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
 Data File : BP004375.D
 Acq On : 19 Dec 2020 01:21
 Operator : CG/JU
 Sample : L5128-12
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB8D6

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 0.1 % 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:\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	3.317	19	22	24	rBV	42457	52582	1.17%	0.090%
2	3.352	24	28	39	rVB	248618	359790	7.98%	0.614%
3	4.116	156	158	163	rVB	41029	47621	1.06%	0.081%
4	4.481	215	220	229	rVB	142200	226388	5.02%	0.387%
5	6.775	607	610	613	rBV4	10422	15873	0.35%	0.027%
6	6.993	641	647	658	rVB2	142572	351677	7.80%	0.601%
7	7.157	669	675	683	rBV	400586	665248	14.75%	1.136%
8	7.357	700	709	717	rVB	478890	791018	17.53%	1.351%
9	7.822	780	788	795	rBV	914921	1489986	33.03%	2.544%
10	7.893	795	800	805	rVB4	32147	54423	1.21%	0.093%
11	8.199	847	852	858	rVB3	23178	45532	1.01%	0.078%
12	8.528	901	908	915	rBV	404290	649051	14.39%	1.108%
13	8.716	937	940	948	rVB5	16352	32807	0.73%	0.056%
14	8.804	948	955	970	rVV	2273284	3704155	82.11%	6.326%
15	8.922	970	975	979	rVV3	33167	60129	1.33%	0.103%
16	8.975	979	984	994	rVV	498750	883539	19.59%	1.509%
17	9.063	994	999	1008	rVV	75420	135822	3.01%	0.232%
18	9.181	1011	1019	1028	rVB9	18693	47485	1.05%	0.081%
19	9.340	1040	1046	1051	rBV5	38391	70767	1.57%	0.121%
20	9.428	1051	1061	1068	rVB2	85004	170960	3.79%	0.292%
21	9.522	1072	1077	1081	rBV7	12449	26728	0.59%	0.046%
22	9.563	1081	1084	1089	rVB7	10367	17279	0.38%	0.030%
23	9.616	1089	1093	1101	rVB8	13481	32758	0.73%	0.056%
24	9.699	1101	1107	1114	rBV	436383	742723	16.46%	1.268%
25	9.834	1123	1130	1134	rBV6	29551	56908	1.26%	0.097%
26	9.881	1134	1138	1144	rVV	35418	63552	1.41%	0.109%
27	10.034	1158	1164	1171	rVB	193393	330310	7.32%	0.564%
28	10.110	1171	1177	1179	rBV4	10269	20421	0.45%	0.035%
29	10.169	1179	1187	1192	rBV	99225	212533	4.71%	0.363%
30	10.240	1192	1199	1218	rVB	755099	1461410	32.40%	2.496%
31	10.393	1218	1225	1230	rBV5	30030	72747	1.61%	0.124%
32	10.528	1243	1248	1256	rVB	43531	83651	1.85%	0.143%
33	10.622	1256	1264	1270	rBV	1262628	2211081	49.01%	3.776%
34	10.675	1270	1273	1276	rVV5	17452	23907	0.53%	0.041%

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35	10.751	1277	1286	1300	rVB	685858	1263031	28.00%	2.157%
36	10.869	1300	1306	1314	rBV8	25436	68724	1.52%	0.117%
37	10.975	1320	1324	1328	rVB6	11688	18697	0.41%	0.032%
38	11.046	1328	1336	1342	rBV9	18198	41196	0.91%	0.070%
39	11.181	1353	1359	1360	rBV6	15212	23524	0.52%	0.040%
40	11.228	1360	1367	1371	rVB8	15398	40772	0.90%	0.070%
41	11.287	1371	1377	1379	rBV5	10719	19085	0.42%	0.033%
42	11.328	1379	1384	1388	rVB6	13124	26641	0.59%	0.045%
43	11.463	1395	1407	1413	rBV5	40526	116258	2.58%	0.199%
44	11.534	1415	1419	1423	rVV5	11441	18236	0.40%	0.031%
45	11.587	1423	1428	1435	rVB6	27366	53589	1.19%	0.092%
46	11.651	1435	1439	1443	rBV6	12363	22483	0.50%	0.038%
47	11.822	1463	1468	1474	rVB2	31147	51729	1.15%	0.088%
48	11.887	1474	1479	1482	rBV4	10069	15901	0.35%	0.027%
49	11.963	1484	1492	1505	rVB	68536	156101	3.46%	0.267%
50	12.069	1505	1510	1513	rBV6	13928	23235	0.52%	0.040%
51	12.122	1513	1519	1528	rBV	1107491	1777743	39.41%	3.036%
52	12.228	1528	1537	1544	rVB	205225	380049	8.42%	0.649%
53	12.340	1550	1556	1560	rBV5	21778	42177	0.93%	0.072%
54	12.504	1580	1584	1588	rVB6	12265	19295	0.43%	0.033%
55	12.557	1589	1593	1595	rBV4	10367	16557	0.37%	0.028%
56	12.622	1600	1604	1608	rVB5	15999	24737	0.55%	0.042%
57	12.716	1617	1620	1625	rVB6	13751	19497	0.43%	0.033%
58	12.840	1635	1641	1649	rBV6	17640	35270	0.78%	0.060%
59	13.181	1694	1699	1706	rVV8	23229	59092	1.31%	0.101%
60	13.251	1707	1711	1717	rVV5	23102	53399	1.18%	0.091%
61	13.316	1717	1722	1728	rVB6	26146	57913	1.28%	0.099%
62	13.375	1728	1732	1736	rBV7	11361	23791	0.53%	0.041%
63	13.504	1750	1754	1760	rBV2	29332	68567	1.52%	0.117%
64	13.557	1760	1763	1774	rVB	45553	84350	1.87%	0.144%
65	13.698	1782	1787	1792	rBV8	16132	37240	0.83%	0.064%
66	13.792	1799	1803	1805	rBV5	8817	13678	0.30%	0.023%
67	13.875	1811	1817	1822	rBV	1334792	1867972	41.41%	3.190%
68	13.922	1822	1825	1831	rVB	132234	174349	3.86%	0.298%
69	14.022	1836	1842	1848	rBV7	34317	75591	1.68%	0.129%
70	14.069	1848	1850	1856	rVB7	14250	21104	0.47%	0.036%
71	14.163	1857	1866	1881	rBV	1509685	2299675	50.98%	3.927%

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72	14.275	1881	1885	1892	rBV	59748	100874	2.24%	0.172%
73	14.392	1900	1905	1910	rBV2	105782	150628	3.34%	0.257%
74	14.469	1912	1918	1925	rVB	1980756	3062064	67.88%	5.229%
75	14.675	1948	1953	1960	rBV	121271	183537	4.07%	0.313%
76	15.216	2041	2045	2050	rBV	39979	74892	1.66%	0.128%
77	15.334	2062	2065	2068	rBV4	12963	16712	0.37%	0.029%
78	15.469	2082	2088	2096	rVB	1899984	2695299	59.75%	4.603%
79	15.592	2103	2109	2124	rVB	377551	593038	13.15%	1.013%
80	15.722	2124	2131	2134	rBV5	70158	143660	3.18%	0.245%
81	15.981	2170	2175	2181	rBV	444073	607647	13.47%	1.038%
82	16.151	2197	2204	2209	rBV2	186027	316627	7.02%	0.541%
83	16.316	2228	2232	2236	rBV3	25343	37975	0.84%	0.065%
84	16.369	2238	2241	2254	rVB5	22919	47821	1.06%	0.082%
85	16.492	2256	2262	2267	rVV	286898	434201	9.62%	0.741%
86	16.545	2267	2271	2279	rVB3	53459	88841	1.97%	0.152%
87	16.763	2305	2308	2313	rBV	55264	72695	1.61%	0.124%
88	17.233	2381	2388	2397	rVB	2473644	3685762	81.70%	6.294%
89	17.328	2398	2404	2414	rVV	2066680	3088721	68.47%	5.275%
90	17.416	2415	2419	2424	rVB	38946	60450	1.34%	0.103%
91	18.128	2536	2540	2548	rBV	104528	189734	4.21%	0.324%
92	18.410	2585	2588	2592	rBV4	38171	51555	1.14%	0.088%
93	18.875	2663	2667	2675	rVB	118825	169412	3.76%	0.289%
94	19.569	2780	2785	2790	rBV	2781284	3830135	84.90%	6.541%
95	19.616	2790	2793	2800	rVB	245430	329141	7.30%	0.562%
96	21.039	3031	3035	3042	rBV2	551388	875256	19.40%	1.495%
97	21.104	3042	3046	3052	rVB	285411	396825	8.80%	0.678%
98	21.321	3078	3083	3091	rBV	3295207	4364281	96.74%	7.453%
99	23.527	3451	3458	3472	rVB	1980448	4057838	89.95%	6.929%
100	23.680	3477	3484	3494	rVB	2305615	4511209	100.00%	7.704%

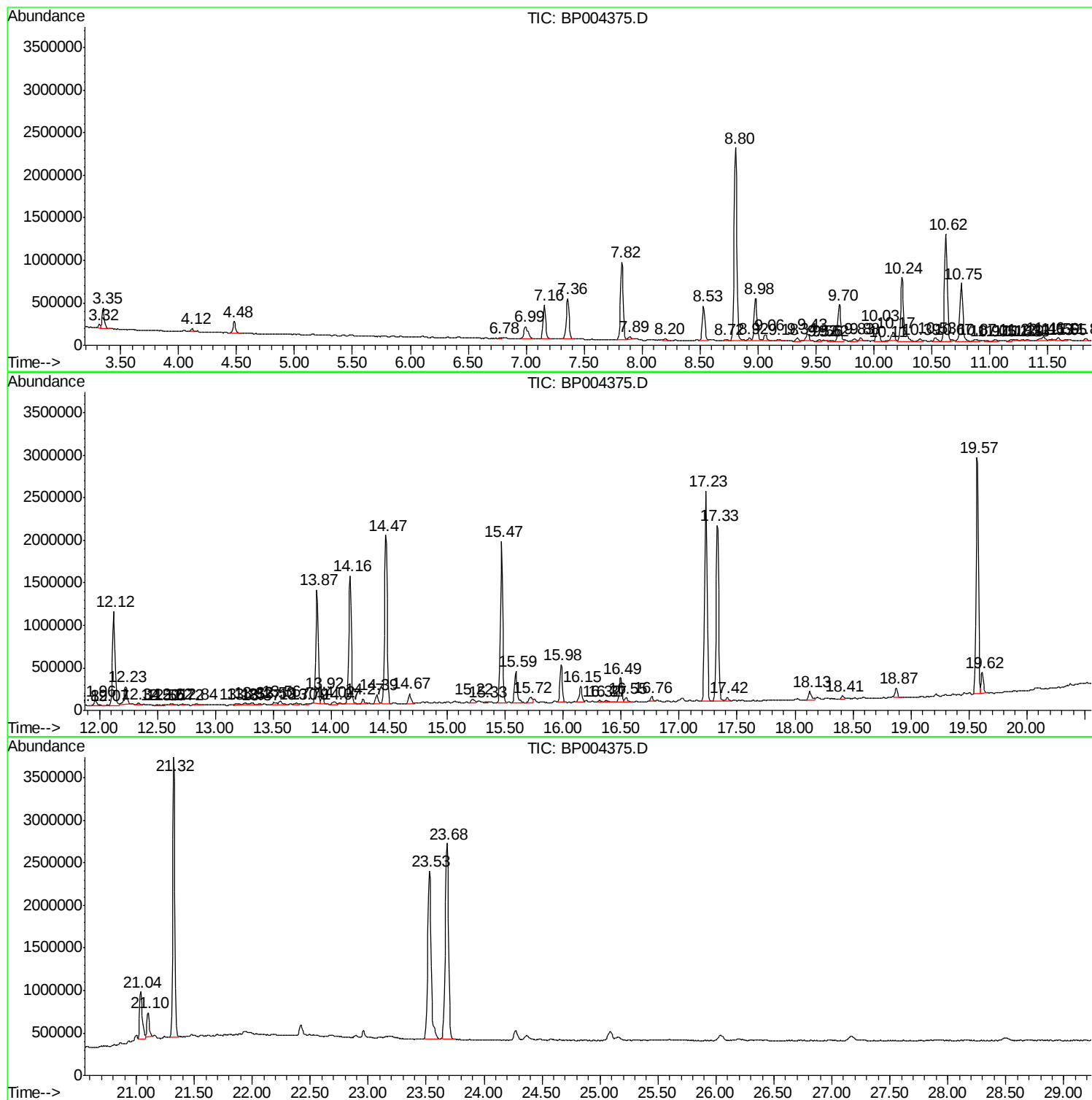
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

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



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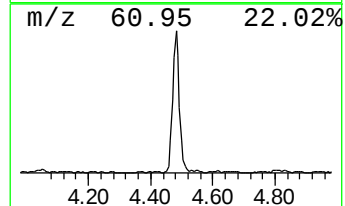
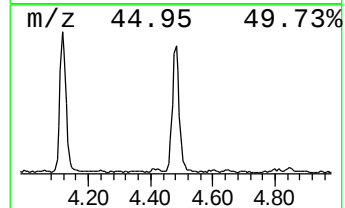
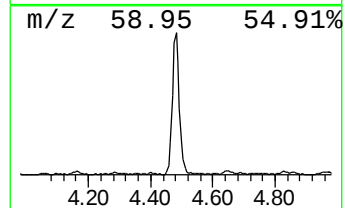
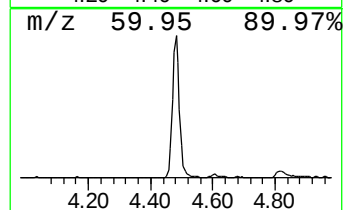
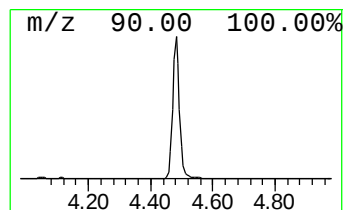
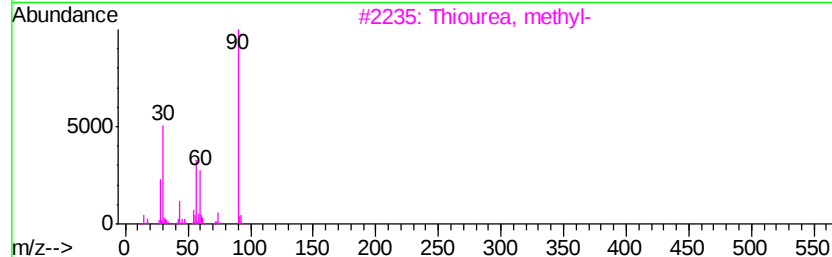
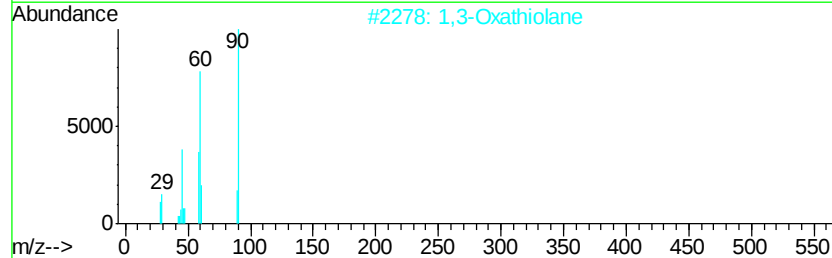
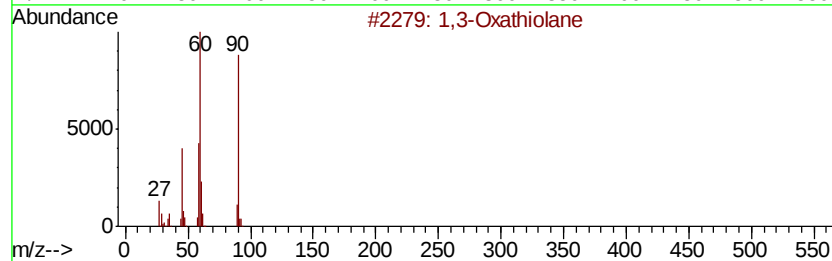
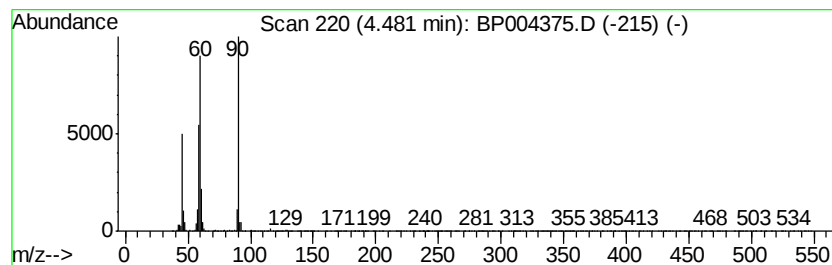
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

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	3.04 ng/ul	226388	1,4-Dichlorobenzene-d4	7.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Oxathiolane	90	C3H6OS	002094-97-5	94
2			1,3-Oxathiolane	90	C3H6OS	002094-97-5	86
3			Thiourea, methyl-	90	C2H6N2S	000598-52-7	50
4			Thiourea, methyl-	90	C2H6N2S	000598-52-7	36
5			Heptanoic acid, 2-hydroxy-, meth...	160	C8H16O3	054340-91-9	36



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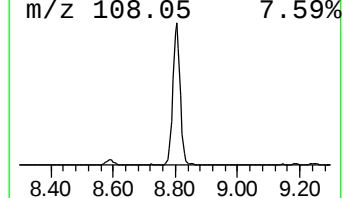
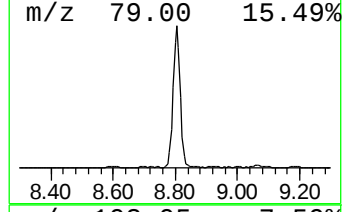
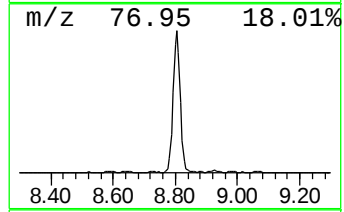
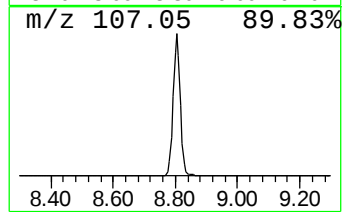
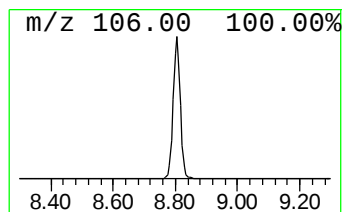
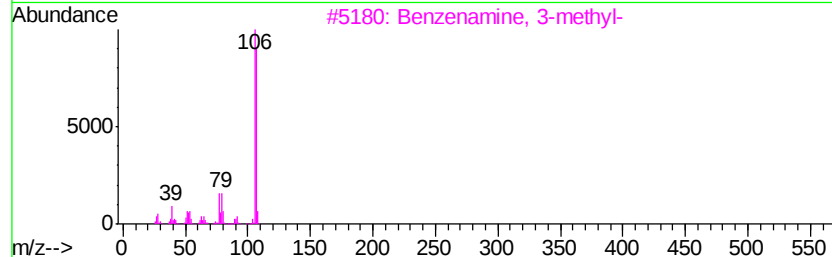
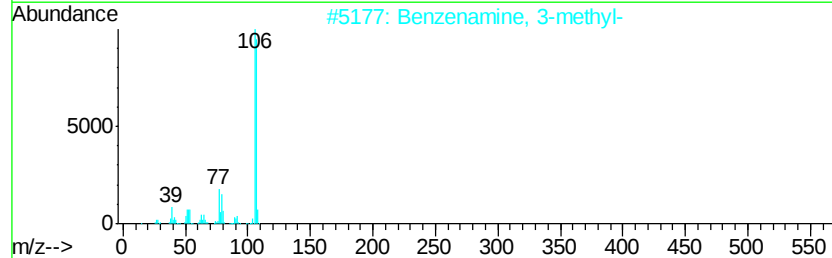
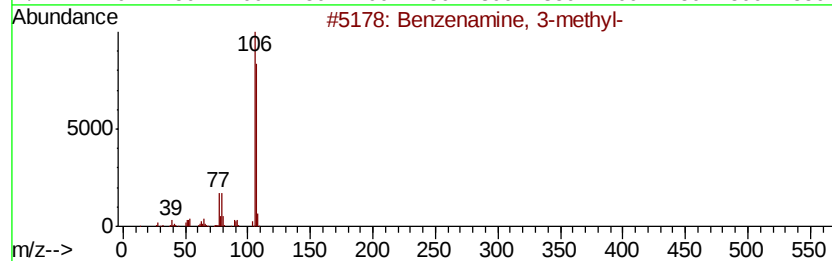
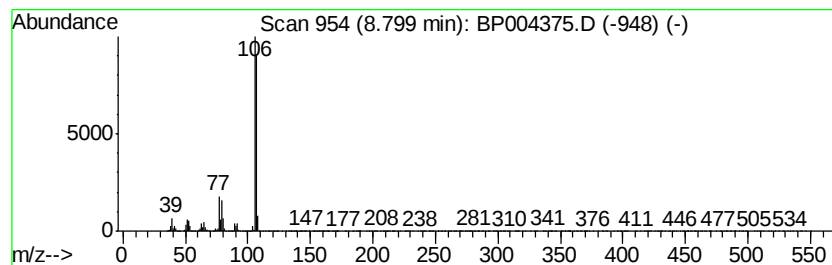
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Benzenamine, 3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.80	49.72 ng/ul	3704160	1,4-Dichlorobenzene-d4	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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



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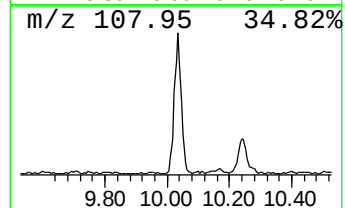
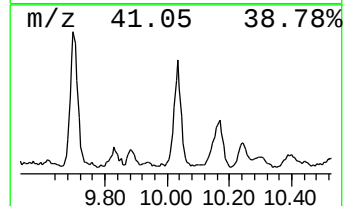
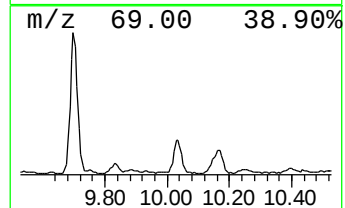
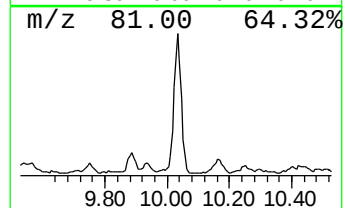
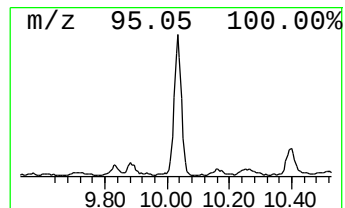
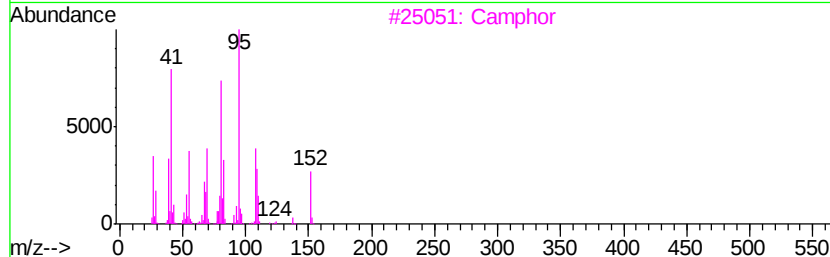
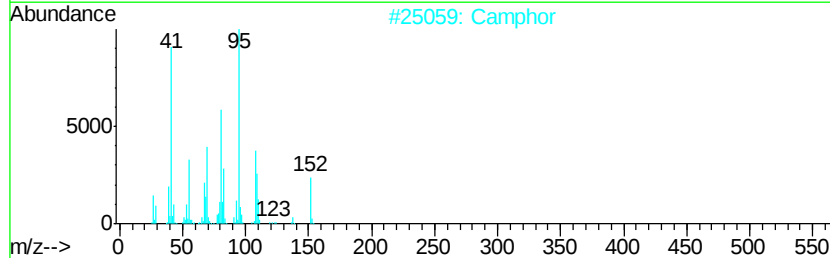
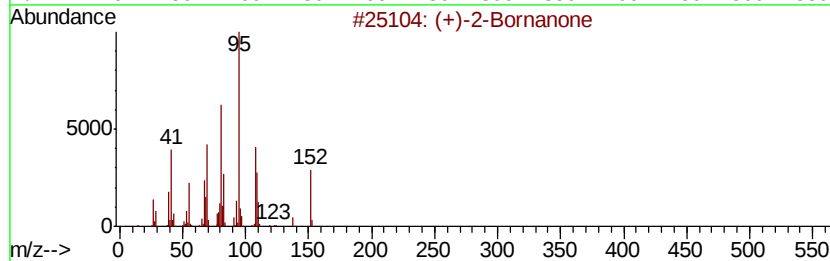
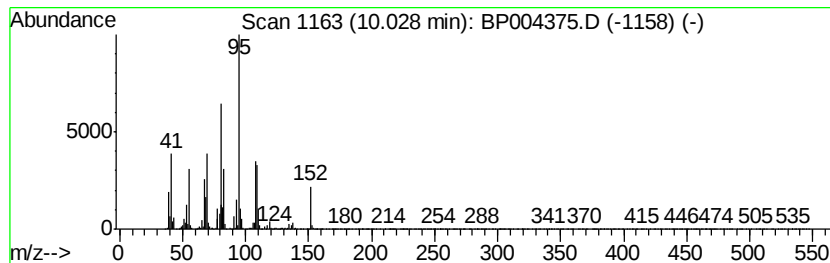
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Peak Number 3 (+)-2-Bornanone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.03	2.99 ng/ul	330310	Naphthalene-d8	10.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	(+)-2-Bornanone	152 C10H16O	000464-49-3	98
2	Camphor	152 C10H16O	000076-22-2	98
3	Camphor	152 C10H16O	000076-22-2	98
4	(+)-2-Bornanone	152 C10H16O	000464-49-3	97
5	(+)-2-Bornanone	152 C10H16O	000464-49-3	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004375.D
Acq On : 19 Dec 2020 01:21
Operator : CG/JU
Sample : L5128-12
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8D6

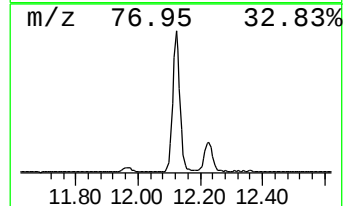
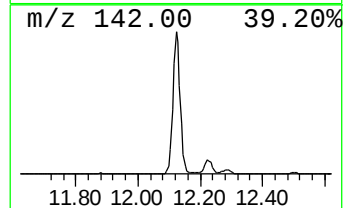
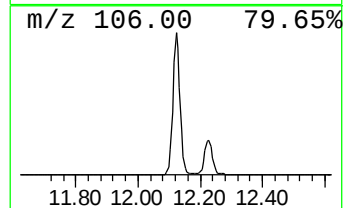
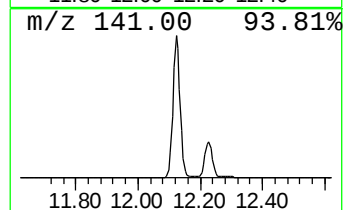
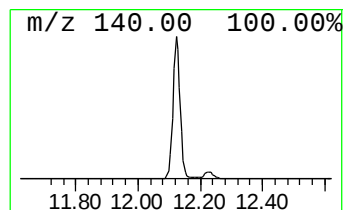
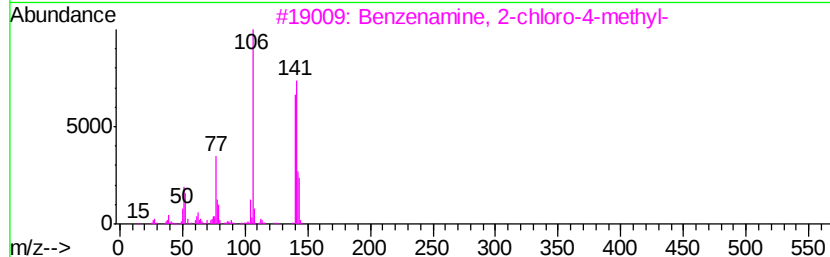
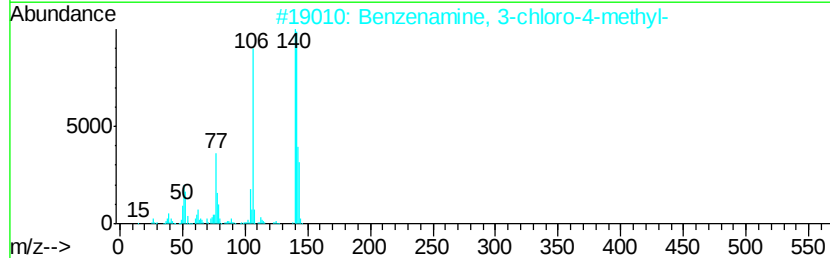
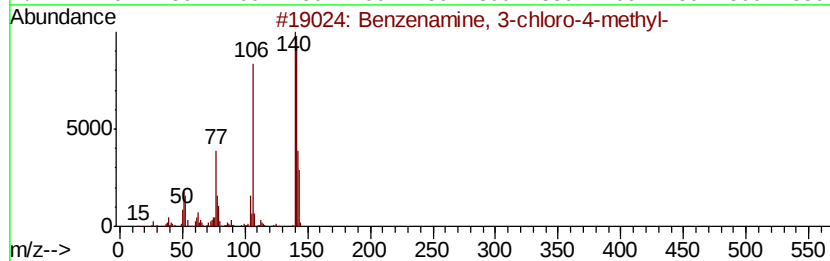
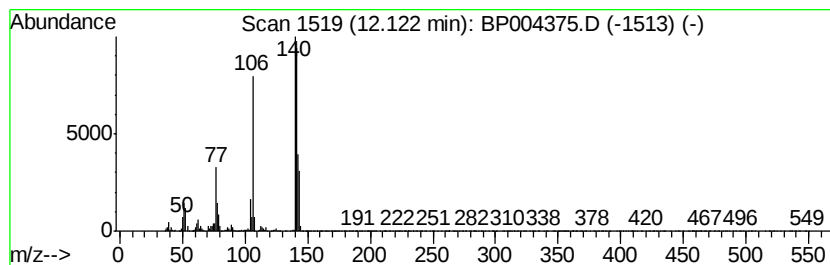
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Quant Title : SVOA CALIBRATION

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	16.08 ng/ul	1777740	Napthalene-d8	10.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004375.D
Acq On : 19 Dec 2020 01:21
Operator : CG/JU
Sample : L5128-12
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8D6

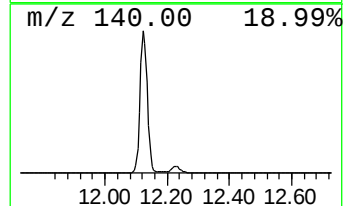
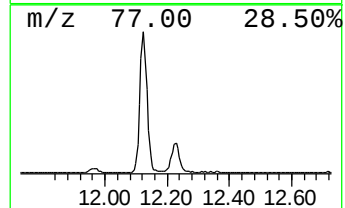
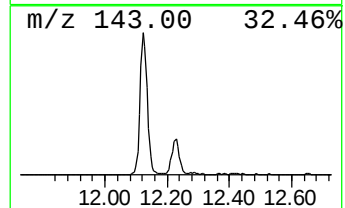
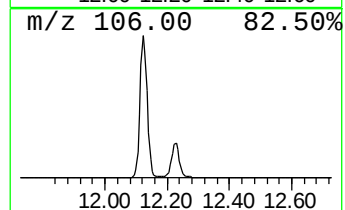
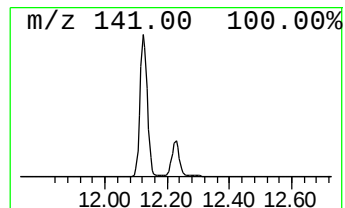
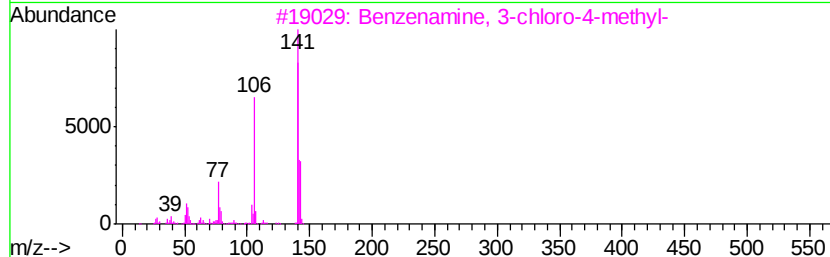
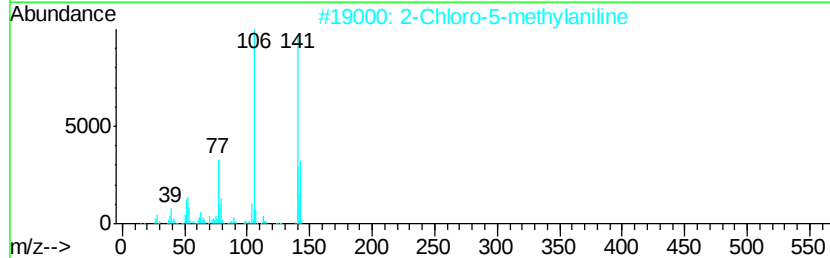
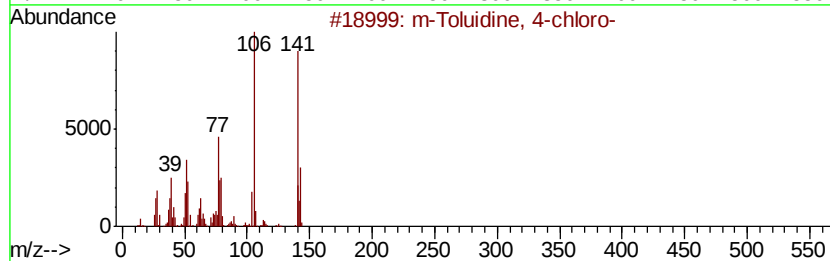
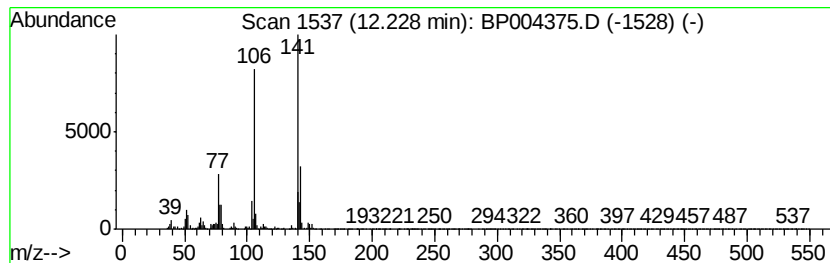
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 m-Toluidine, 4-chloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.23	3.44 ng/ul	380049	Napthalene-d8	10.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	m-Toluidine, 4-chloro-	141	C7H8ClN	007149-75-9	93
2		2-Chloro-5-methylaniline	141	C7H8ClN	000095-81-8	93
3		Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	90
4		Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	87
5		Benzenamine, 4-chloro-2-methyl-	141	C7H8ClN	000095-69-2	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004375.D
Acq On : 19 Dec 2020 01:21
Operator : CG/JU
Sample : L5128-12
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8D6

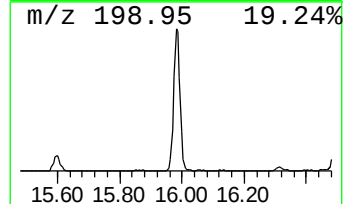
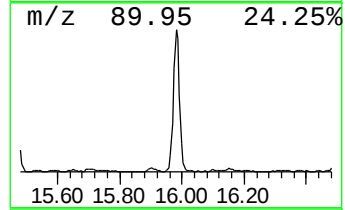
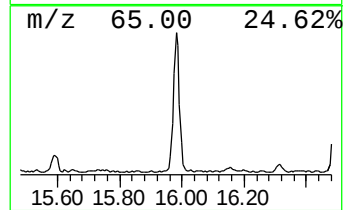
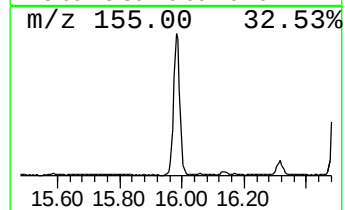
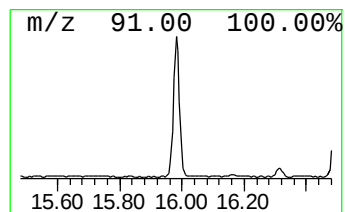
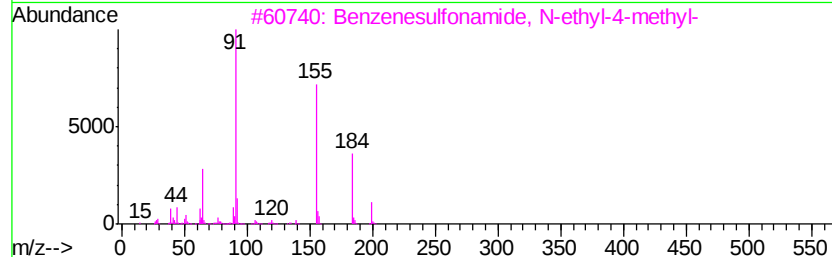
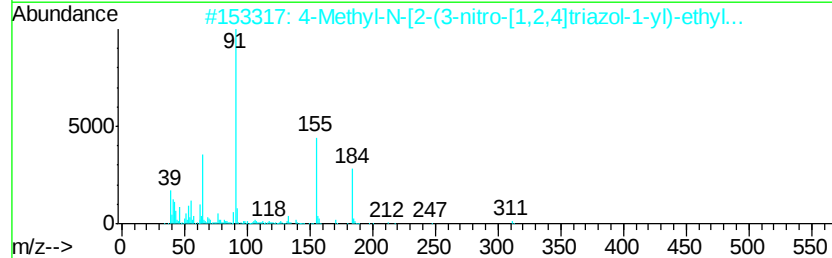
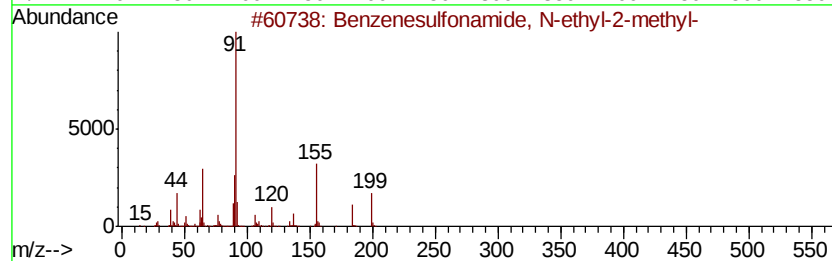
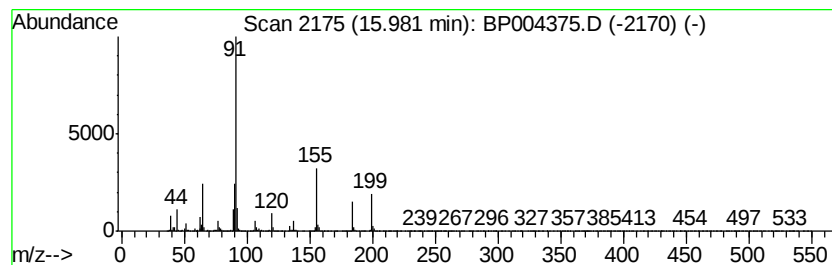
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzenesulfonamide, N-ethyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.98	3.30 ng/ul	607647	Phenanthrene-d10	17.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	60
3		Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	53
4		Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	52
5		Benzenesulfonamide, N,N'-(dithio...	460	C18H24N2O4S4	023516-74-7	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004375.D
Acq On : 19 Dec 2020 01:21
Operator : CG/JU
Sample : L5128-12
Misc :
ALS Vial : 22 Sample Multiplier: 1

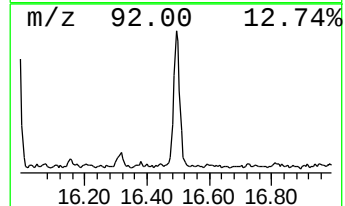
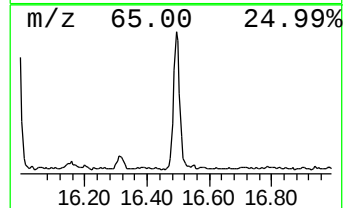
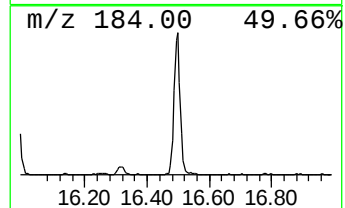
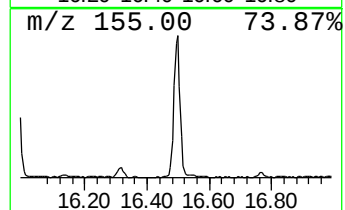
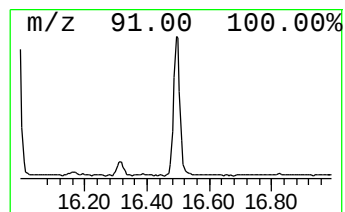
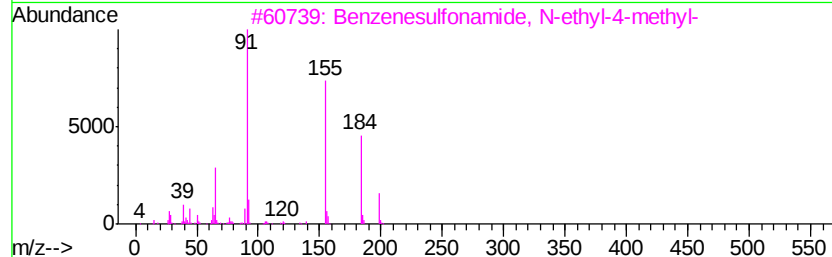
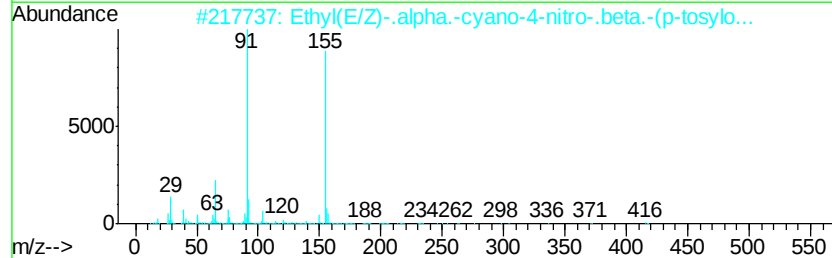
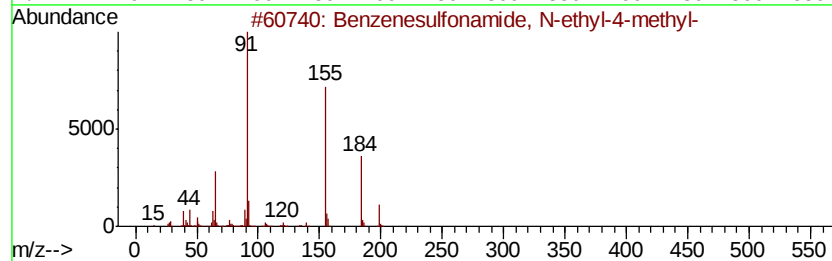
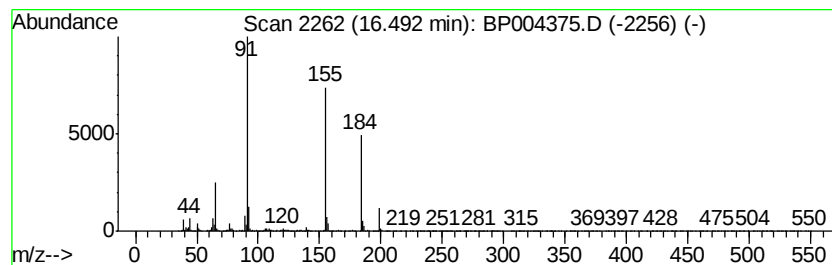
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzenesulfonamide, N-ethyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.49	2.36 ng/ul	434201	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		Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	93
3		Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	91
4		4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	90
5		(Toluene-4-sulfonylamino)-acetic...	283	C13H17NO4S	1000190-48-0	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004375.D
Acq On : 19 Dec 2020 01:21
Operator : CG/JU
Sample : L5128-12
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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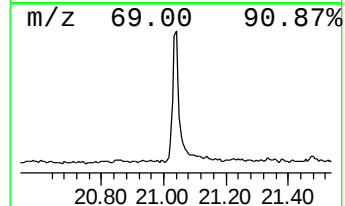
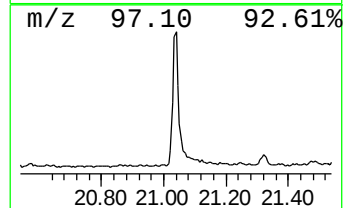
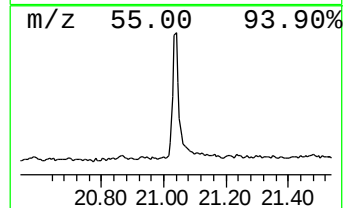
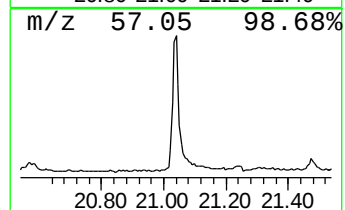
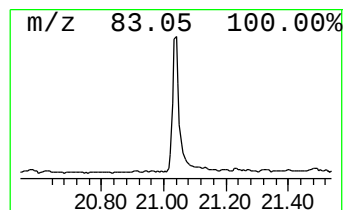
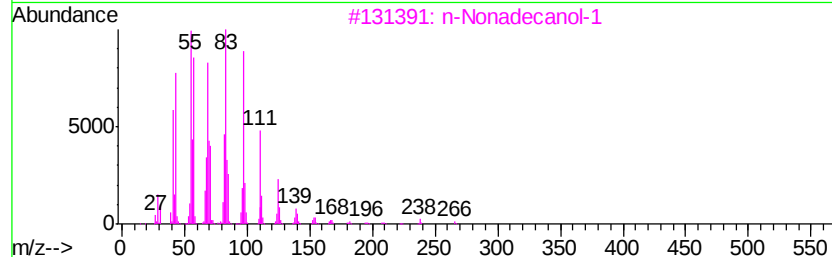
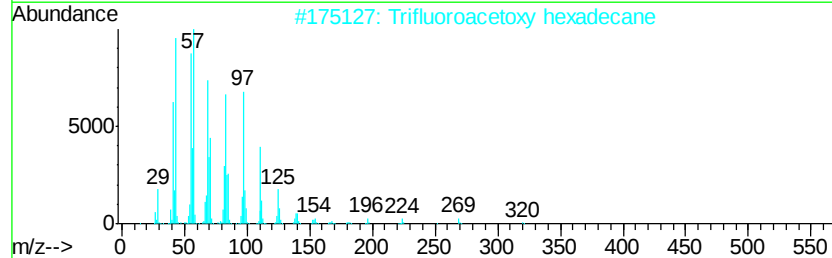
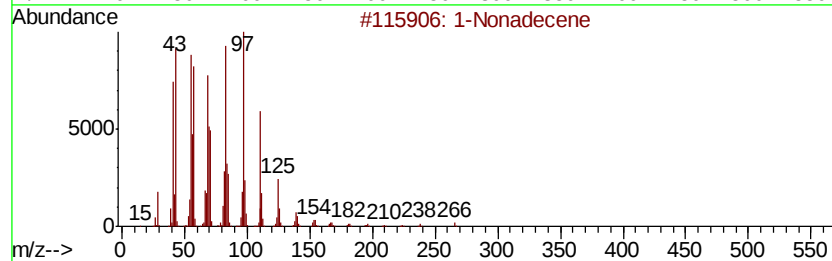
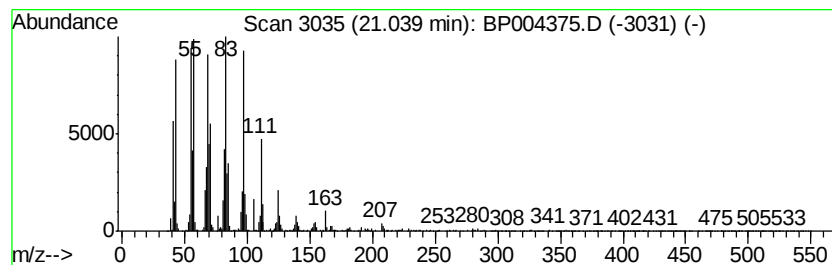
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.04	4.01 ng/ul	875256	Chrysene-d12	21.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Nonadecene	266	C19H38	018435-45-5	94
2			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
3			n-Nonadecanol-1	284	C19H40O	001454-84-8	94
4			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5			1-Heneicosanol	312	C21H44O	015594-90-8	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP121820\
 Data File : BP004375.D
 Acq On : 19 Dec 2020 01:21
 Operator : CG/JU
 Sample : L5128-12
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,3-Oxathiolane	4.48	3.0	ng/ul	226388	1	7.82	1489990	20.0
Benzenamine, 3-me...	8.80	49.7	ng/ul	3704160	1	7.82	1489990	20.0
(+)-2-Bornanone	10.03	3.0	ng/ul	330310	2	10.62	2211080	20.0
Benzenamine, 3-ch...	12.12	16.1	ng/ul	1777740	2	10.62	2211080	20.0
m-Toluidine, 4-ch...	12.23	3.4	ng/ul	380049	2	10.62	2211080	20.0
Benzenesulfonamid...	15.98	3.3	ng/ul	607647	4	17.23	3685760	20.0
Benzenesulfonamid...	16.49	2.4	ng/ul	434201	4	17.23	3685760	20.0
(DEL) Alkane: Str...	21.04	4.0	ng/ul	875256	5	21.32	4364280	20.0