

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
 Data File : BP014483.D
 Acq On : 03 Apr 2023 12:59
 Operator : CG/JU
 Sample : 02093-05DL 10X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB973DL

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\SFAM-EPA-BP032823.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014483.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.387	32	34	35	rBV2	4287	3354	0.11%	0.018%
2	3.422	35	40	54	rVB	165669	242890	8.18%	1.273%
3	3.669	78	82	88	rBV	18395	25065	0.84%	0.131%
4	3.875	113	117	121	rBV4	10258	19919	0.67%	0.104%
5	4.140	158	162	168	rBV	11820	16330	0.55%	0.086%
6	4.216	171	175	184	rVB	31308	45733	1.54%	0.240%
7	4.599	232	240	248	rVV	26968	50411	1.70%	0.264%
8	5.781	437	441	442	rBV4	2496	3010	0.10%	0.016%
9	6.275	521	525	533	rVB	7191	11826	0.40%	0.062%
10	6.510	560	565	570	rBV5	5465	8854	0.30%	0.046%
11	6.710	595	599	607	rVB4	13791	25169	0.85%	0.132%
12	7.058	653	658	660	rBV6	3075	4709	0.16%	0.025%
13	7.116	660	668	670	rVV6	3741	9388	0.32%	0.049%
14	7.199	677	682	691	rBV2	15763	43054	1.45%	0.226%
15	7.334	699	705	715	rBV	69152	128470	4.32%	0.674%
16	7.546	735	741	753	rBV2	66140	129494	4.36%	0.679%
17	7.622	753	754	759	rVB5	3661	4378	0.15%	0.023%
18	8.005	813	819	830	rBV	663896	1109243	37.34%	5.815%
19	8.093	830	834	841	rVB9	8066	16068	0.54%	0.084%
20	8.746	938	945	959	rBV2	49308	125188	4.21%	0.656%
21	8.840	959	961	965	rVV5	4618	6697	0.23%	0.035%
22	8.893	968	970	973	rVB4	3266	3074	0.10%	0.016%
23	9.122	1005	1009	1014	rVV8	3922	7040	0.24%	0.037%
24	9.193	1014	1021	1032	rBV	80172	152763	5.14%	0.801%
25	9.387	1050	1054	1058	rVB7	2090	3358	0.11%	0.018%
26	9.446	1060	1064	1069	rVB8	2334	3679	0.12%	0.019%
27	9.534	1073	1079	1087	rBV8	11347	27302	0.92%	0.143%
28	9.634	1092	1096	1102	rVB9	4127	5522	0.19%	0.029%
29	9.740	1109	1114	1116	rBV6	3503	5693	0.19%	0.030%
30	9.757	1116	1117	1122	rVV5	4183	4657	0.16%	0.024%
31	9.840	1124	1131	1137	rVB5	4545	11070	0.37%	0.058%
32	9.916	1137	1144	1155	rBV	65668	131510	4.43%	0.689%
33	10.181	1185	1189	1190	rBV4	2358	3178	0.11%	0.017%
34	10.381	1216	1223	1229	rVB2	24612	48734	1.64%	0.255%
35	10.481	1232	1240	1251	rBV	72986	199773	6.73%	1.047%
36	10.834	1291	1300	1313	rBV	1014701	1775053	59.76%	9.306%

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
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37	10.993	1317	1327	1338	rBV	80410	194285	6.54%	1.019%
38	11.422	1396	1400	1406	rVB8	9461	16409	0.55%	0.086%
39	11.534	1417	1419	1425	rVB7	3179	4394	0.15%	0.023%
40	11.657	1433	1440	1442	rBV8	6965	14149	0.48%	0.074%
41	11.816	1464	1467	1468	rBV3	3529	3749	0.13%	0.020%
42	12.040	1500	1505	1512	rVB6	13761	27611	0.93%	0.145%
43	12.145	1519	1523	1526	rBV6	2858	4611	0.16%	0.024%
44	12.187	1526	1530	1539	rVB6	10394	20785	0.70%	0.109%
45	12.357	1549	1559	1569	rBV4	23287	69536	2.34%	0.365%
46	12.475	1574	1579	1588	rVV4	10581	32184	1.08%	0.169%
47	12.551	1588	1592	1599	rVB9	4490	11397	0.38%	0.060%
48	12.651	1604	1609	1615	rVB3	33575	56423	1.90%	0.296%
49	12.928	1653	1656	1660	rBV6	4724	5040	0.17%	0.026%
50	13.051	1672	1677	1680	rBV7	5542	8832	0.30%	0.046%
51	13.216	1701	1705	1710	rBV7	9347	15791	0.53%	0.083%
52	13.463	1743	1747	1753	rBV9	4160	9017	0.30%	0.047%
53	13.545	1756	1761	1766	rBV3	16733	32685	1.10%	0.171%
54	13.710	1785	1789	1794	rBV4	11460	19946	0.67%	0.105%
55	13.892	1816	1820	1827	rBV2	22563	44386	1.49%	0.233%
56	14.069	1844	1850	1862	rBV	253879	371096	12.49%	1.946%
57	14.175	1864	1868	1875	rBV2	21239	37797	1.27%	0.198%
58	14.357	1894	1899	1915	rBV2	253733	411144	13.84%	2.156%
59	14.469	1915	1918	1929	rVB3	21123	40501	1.36%	0.212%
60	14.581	1933	1937	1942	rBV3	59247	83611	2.81%	0.438%
61	14.663	1944	1951	1964	rVV	1647490	2472335	83.23%	12.962%
62	15.034	2010	2014	2019	rVB4	20322	27816	0.94%	0.146%
63	15.404	2073	2077	2082	rBV	25667	43517	1.46%	0.228%
64	15.657	2115	2120	2131	rVB	356474	525464	17.69%	2.755%
65	15.792	2139	2143	2150	rVB3	57079	97852	3.29%	0.513%
66	16.181	2204	2209	2218	rVB	238135	330998	11.14%	1.735%
67	16.357	2236	2239	2243	rBV6	19075	30100	1.01%	0.158%
68	16.698	2292	2297	2302	rBV	123148	190176	6.40%	0.997%
69	16.951	2338	2340	2347	rVB6	15556	19790	0.67%	0.104%
70	17.428	2414	2421	2429	rBV	2002589	2970475	100.00%	15.573%
71	17.528	2432	2438	2446	rVV	368482	562572	18.94%	2.949%
72	18.292	2565	2568	2577	rBV10	29849	44271	1.49%	0.232%
73	19.051	2695	2697	2701	rVB5	33942	35882	1.21%	0.188%
74	19.757	2812	2817	2822	rVB	455861	600225	20.21%	3.147%
75	21.516	3111	3116	3123	rBV	1852996	2468249	83.09%	12.940%
76	23.868	3513	3516	3523	rVB2	269704	451752	15.21%	2.368%

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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\SFAM-EPA-BP032823.M
Title : SVOA CALIBRATION

77 24.033 3538 3544 3555 rVB2 1066263 2255599 75.93% 11.825%

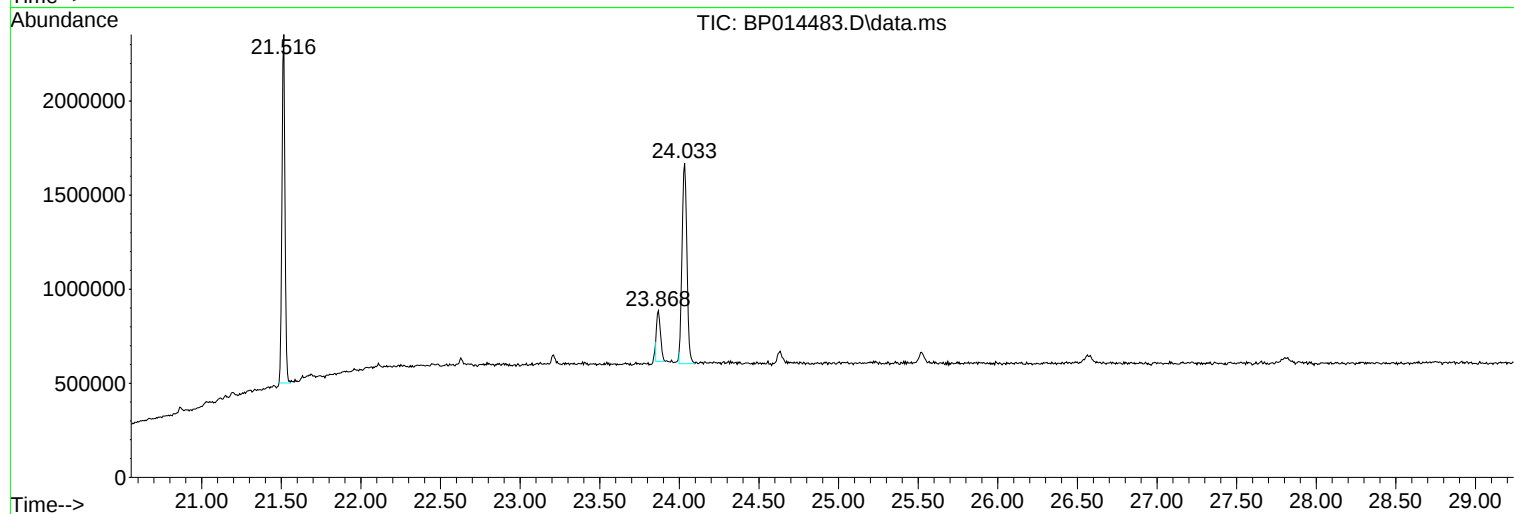
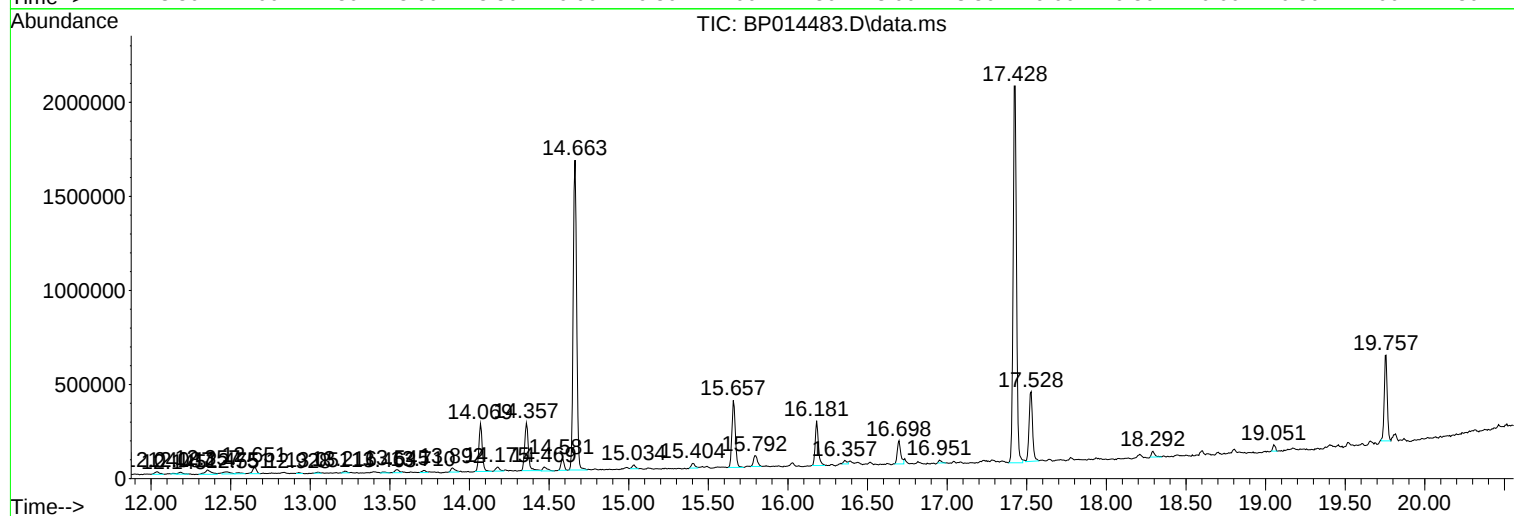
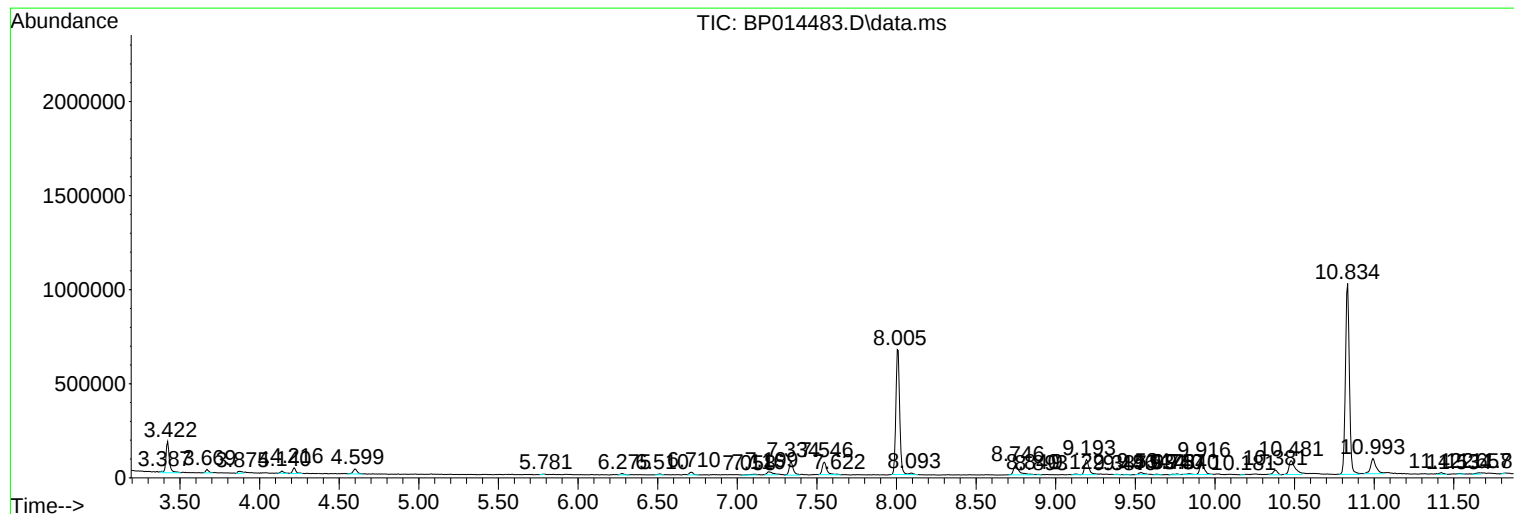
Sum of corrected areas: 19074108

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BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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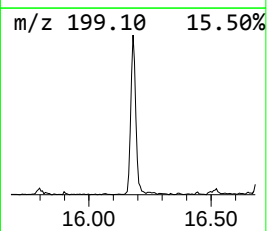
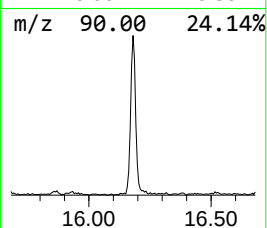
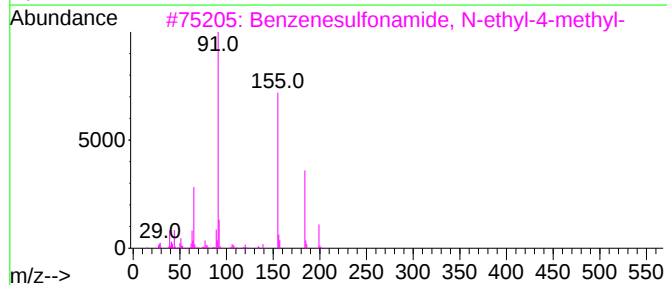
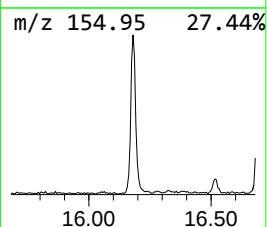
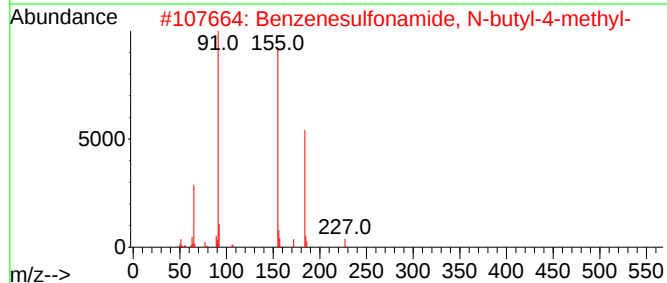
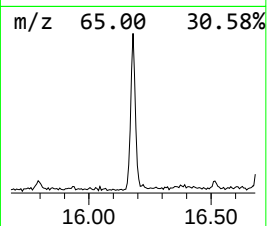
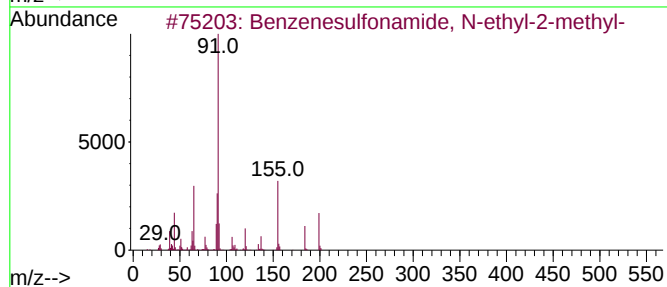
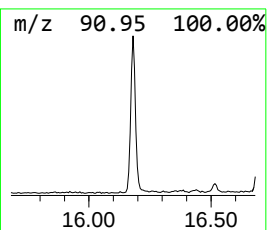
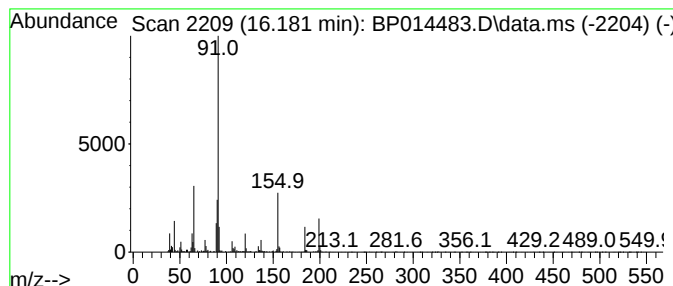
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzenesulfonamide, N-ethyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.181	2.23 ng/ul	330998	Phenanthrene-d10	17.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2			Benzenesulfonamide, N-butyl-4-me...	227	C11H17NO2S	001907-65-9	62
3			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	59
4			Carbamic acid, methyl[(4-methylp...	243	C10H13NO4S	032258-50-7	59
5			p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	58



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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzenesulfonam...	16.181	2.2	ng/u1	330998	4	17.428	2970480	20.0