

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122023\
 Data File : BP018991.D
 Acq On : 21 Dec 2023 09:00
 Operator : MA/JU
 Sample : 05894-18
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH3N3

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\SFAM-EPA-BP120123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018991.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.252	25	28	34	rVB	23062	27453	0.31%	0.056%
2	3.675	97	100	113	rBV	68771	106746	1.21%	0.219%
3	3.899	135	138	144	rVB3	9318	14302	0.16%	0.029%
4	3.999	151	155	159	rVB	9809	13095	0.15%	0.027%
5	5.128	343	347	352	rBV2	8359	11067	0.12%	0.023%
6	6.999	659	665	671	rVB	93404	149674	1.69%	0.308%
7	7.158	686	692	705	rVB	477037	724418	8.18%	1.489%
8	7.352	716	725	735	rBV	437772	679348	7.67%	1.397%
9	7.822	798	805	816	rBV	447120	692347	7.82%	1.423%
10	8.534	918	926	940	rBV	292089	493825	5.58%	1.015%
11	8.981	994	1002	1010	rBV	545306	889696	10.05%	1.829%
12	9.705	1117	1125	1133	rBV	428395	699550	7.90%	1.438%
13	10.240	1207	1216	1233	rBV	610778	1072153	12.11%	2.204%
14	10.616	1273	1280	1288	rBV	585063	943065	10.65%	1.939%
15	10.757	1297	1304	1321	rBV	429062	770736	8.70%	1.584%
16	12.210	1542	1551	1559	rVB4	12864	26485	0.30%	0.054%
17	13.357	1741	1746	1752	rVB4	8344	13065	0.15%	0.027%
18	13.875	1827	1834	1843	rBV	1455670	2074375	23.43%	4.264%
19	14.151	1873	1881	1897	rBV	1508727	2346306	26.50%	4.823%
20	14.457	1926	1933	1942	rBV	930545	1389360	15.69%	2.856%
21	14.681	1964	1971	1989	rBV2	60946	149379	1.69%	0.307%
22	14.881	2000	2005	2011	rVB9	4565	10020	0.11%	0.021%
23	15.451	2093	2102	2117	rBV	2169190	3224179	36.41%	6.628%
24	15.575	2117	2123	2132	rBV	696850	949766	10.73%	1.952%
25	15.687	2139	2142	2148	rVB2	10621	16363	0.18%	0.034%
26	16.239	2231	2236	2243	rVB4	9194	15094	0.17%	0.031%
27	17.034	2361	2371	2376	rBV	5867845	8855001	100.00%	18.203%
28	17.163	2389	2393	2395	rBV5	29453	38484	0.43%	0.079%
29	17.204	2395	2400	2411	rVV	1297953	1845947	20.85%	3.795%
30	17.304	2411	2417	2433	rVB	2847489	3984623	45.00%	8.191%
31	18.098	2547	2552	2559	rBV3	40110	63237	0.71%	0.130%
32	18.169	2562	2564	2571	rVB3	12968	17216	0.19%	0.035%
33	19.033	2706	2711	2716	rBV	74491	90070	1.02%	0.185%
34	19.122	2722	2726	2733	rVB2	43224	59042	0.67%	0.121%
35	19.192	2734	2738	2741	rBV	35503	44502	0.50%	0.091%
36	19.333	2758	2762	2769	rBV3	35532	49449	0.56%	0.102%

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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-BP120123.MA.M
Title : SVOA CALIBRATION

37	19.545	2792	2798	2808	rBV	3999085	5183001	58.53%	10.655%
38	21.298	3091	3096	3103	rBV	1986895	2392250	27.02%	4.918%
39	21.392	3109	3112	3118	rVB	159021	169584	1.92%	0.349%
40	23.504	3464	3471	3485	rVB2	2893435	5825495	65.79%	11.975%
41	23.657	3491	3497	3506	rVB	1247756	2526341	28.53%	5.193%

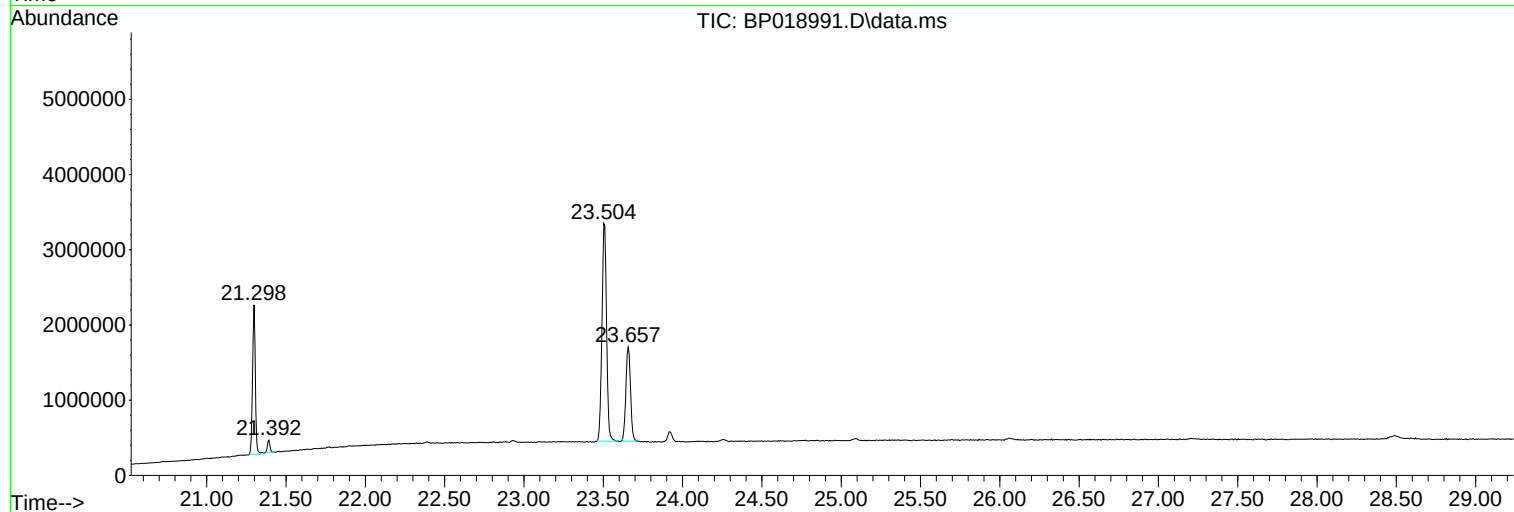
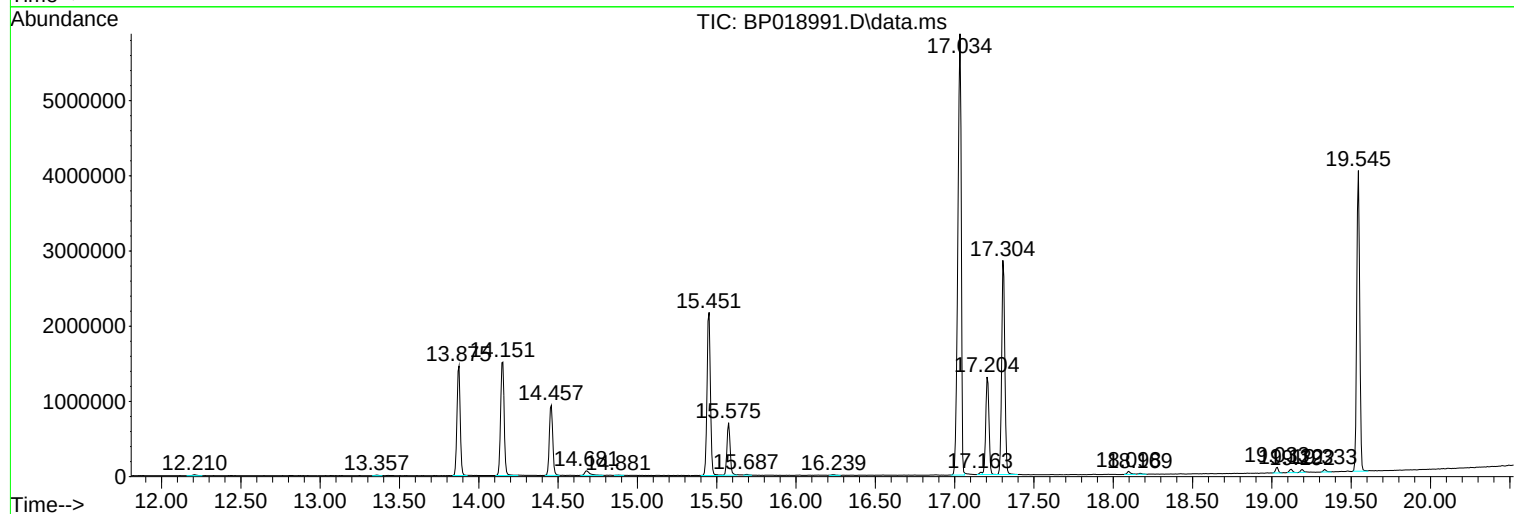
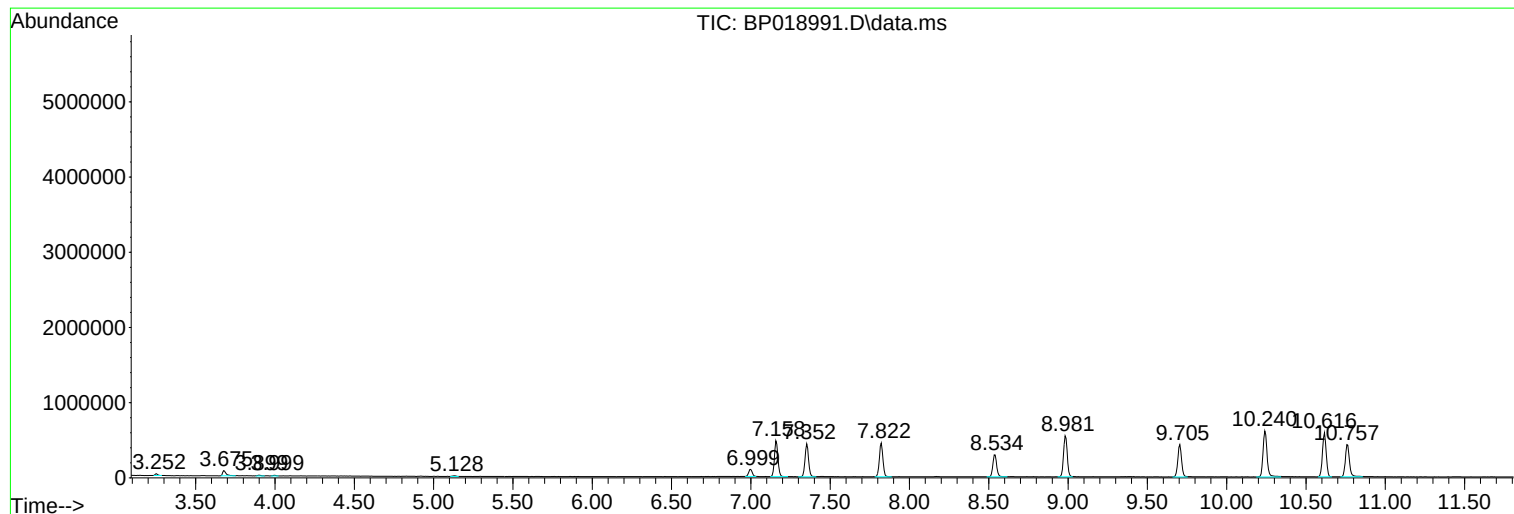
Sum of corrected areas: 48646109

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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



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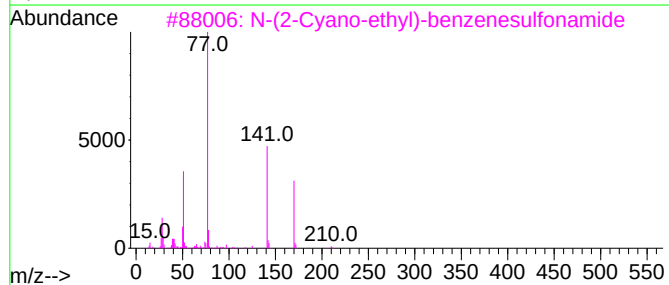
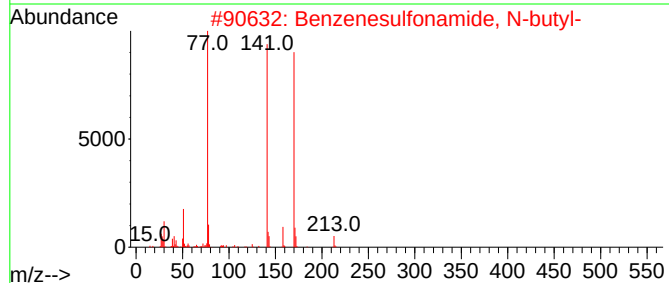
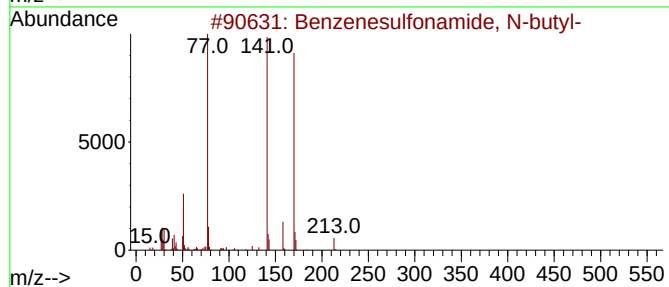
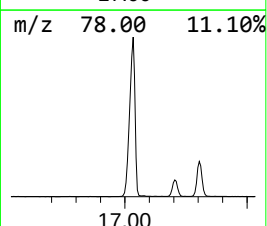
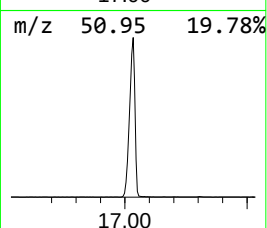
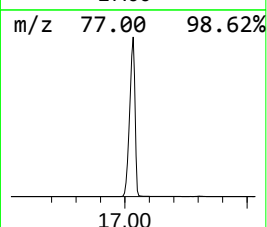
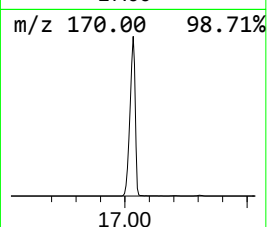
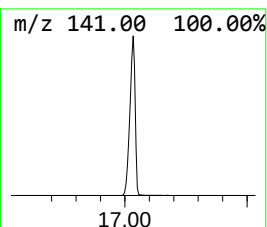
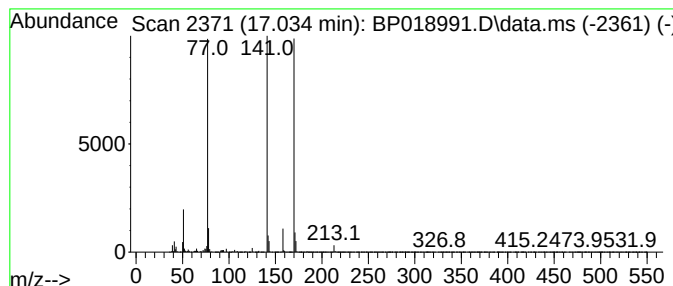
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzenesulfonamide, N-butyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.034	95.94 ng/ul	8855000	Phenanthrene-d10	17.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	90
2			Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	90
3			N-(2-Cyano-ethyl)-benzenesulfona...	210	C9H10N2O2S	002619-21-8	64
4			2-propenoic acid, 2-methyl-, 2-[...	269	C12H15NO4S	1000401-71-8	64
5			N-Benzenesulfonyloxy-2,2-bis(tri...	335	C10H7F6NO3S	055734-41-3	50



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzenesulfonam...	17.034	95.9	ng/ul	8855000	4	17.204	1845950	20.0