

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122123\
 Data File : BM043535.D
 Acq On : 22 Dec 2023 12:12
 Operator : MA/JU
 Sample : 05927-02
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BH3J3

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_M\Methods\SFAM-EPA-BM121523.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM043535.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.369	28	31	37	rVB	45574	64504	1.20%	0.118%
2	3.799	101	104	119	rBV	202573	357302	6.67%	0.655%
3	4.022	139	142	147	rVB2	17035	22580	0.42%	0.041%
4	4.122	157	159	164	rVB	18167	21327	0.40%	0.039%
5	4.575	233	236	240	rVB4	11423	15224	0.28%	0.028%
6	4.952	298	300	304	rVB4	4902	6541	0.12%	0.012%
7	5.052	315	317	321	rVB5	7211	9097	0.17%	0.017%
8	6.187	508	510	515	rBV6	3560	5909	0.11%	0.011%
9	7.128	663	670	679	rVB	221442	355231	6.64%	0.652%
10	7.310	694	701	712	rBV	778725	1270303	23.73%	2.330%
11	7.493	725	732	742	rBV	752345	1197927	22.38%	2.198%
12	7.581	744	747	750	rVB5	4878	6265	0.12%	0.011%
13	7.969	806	813	824	rBV	794582	1265623	23.64%	2.322%
14	8.681	926	934	942	rBV	558996	940223	17.56%	1.725%
15	9.139	1004	1012	1024	rBV	927733	1542108	28.81%	2.829%
16	9.863	1126	1135	1145	rBV	658490	1115902	20.84%	2.047%
17	10.392	1217	1225	1243	rBV	907675	1608660	30.05%	2.951%
18	10.775	1282	1290	1300	rVB	1061569	1794609	33.52%	3.292%
19	10.922	1307	1315	1332	rBV	808999	1450648	27.10%	2.661%
20	12.016	1500	1501	1506	rVB5	5446	7086	0.13%	0.013%
21	12.369	1555	1561	1568	rVB3	15623	28385	0.53%	0.052%
22	13.504	1746	1754	1759	rBV3	22386	42605	0.80%	0.078%
23	13.586	1763	1768	1771	rVB7	4293	7632	0.14%	0.014%
24	14.022	1834	1842	1852	rBV	1809406	2532984	47.31%	4.647%
25	14.298	1882	1889	1906	rBV	2109582	3014974	56.32%	5.531%
26	14.604	1933	1941	1951	rBV	1520075	2213780	41.35%	4.061%
27	14.810	1969	1976	1992	rBV	144097	305557	5.71%	0.561%
28	15.027	2009	2013	2018	rVB8	7026	13734	0.26%	0.025%
29	15.263	2050	2053	2057	rBV6	4003	5990	0.11%	0.011%
30	15.598	2100	2110	2123	rBV	2657850	3856313	72.03%	7.074%
31	15.716	2123	2130	2141	rBV	864944	1214009	22.68%	2.227%
32	15.839	2147	2151	2154	rVB5	10974	15629	0.29%	0.029%
33	16.380	2239	2243	2249	rVB7	9353	17114	0.32%	0.031%
34	16.757	2304	2307	2312	rBV6	4253	7764	0.15%	0.014%
35	17.162	2369	2376	2395	rBV	3547923	4890517	91.35%	8.971%
36	17.351	2401	2408	2418	rBV	1786963	2551357	47.66%	4.680%

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37	17.445	2418	2424	2439	rVV2	3242175	4636789	86.61%	8.506%
38	18.033	2521	2524	2526	rVB3	5065	5458	0.10%	0.010%
39	18.133	2537	2541	2544	rVB6	4094	6925	0.13%	0.013%
40	18.245	2554	2560	2564	rBV2	57577	84301	1.57%	0.155%
41	18.333	2571	2575	2577	rVB3	7740	9023	0.17%	0.017%
42	18.827	2656	2659	2662	rBV5	4201	5852	0.11%	0.011%
43	19.209	2720	2724	2728	rVB5	20453	28394	0.53%	0.052%
44	19.298	2735	2739	2744	rBV3	33159	50057	0.94%	0.092%
45	19.374	2747	2752	2755	rBV3	34959	53464	1.00%	0.098%
46	19.521	2773	2777	2782	rBV2	35438	50955	0.95%	0.093%
47	19.674	2800	2803	2807	rBV4	12702	17590	0.33%	0.032%
48	19.733	2807	2813	2825	rVV	3685302	5117501	95.59%	9.388%
49	21.527	3112	3118	3128	rBV	1929532	2555752	47.74%	4.688%
50	23.786	3494	3502	3513	rBV	2902192	5353493	100.00%	9.821%
51	23.939	3522	3528	3544	rVB	1356711	2762091	51.59%	5.067%

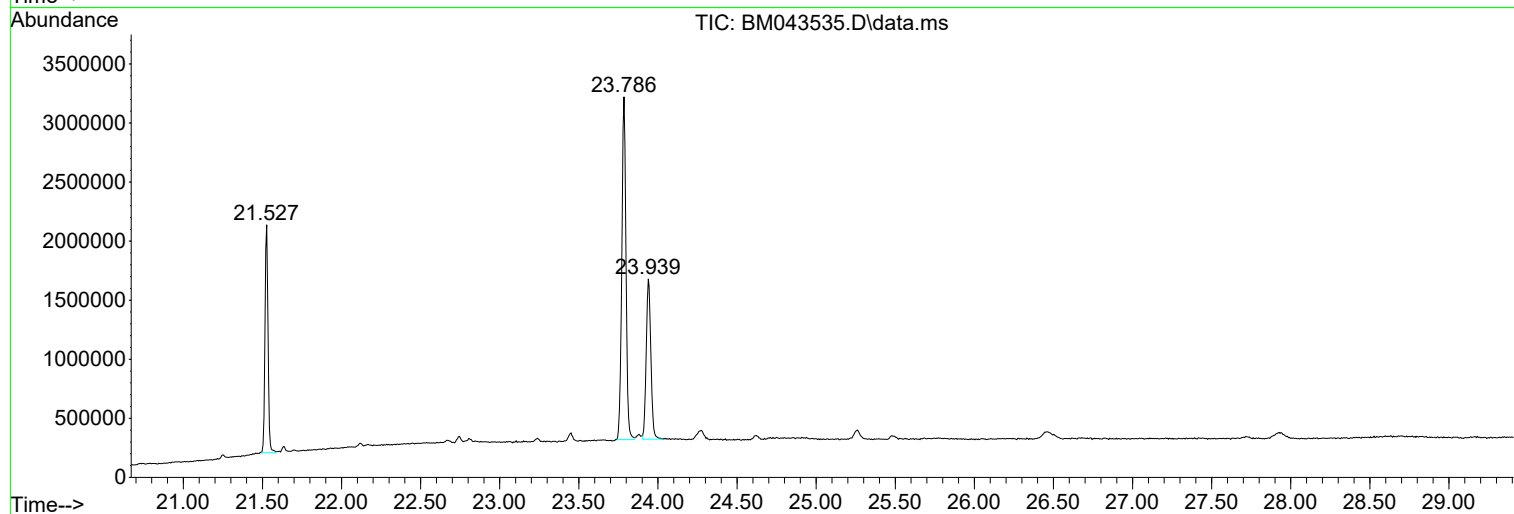
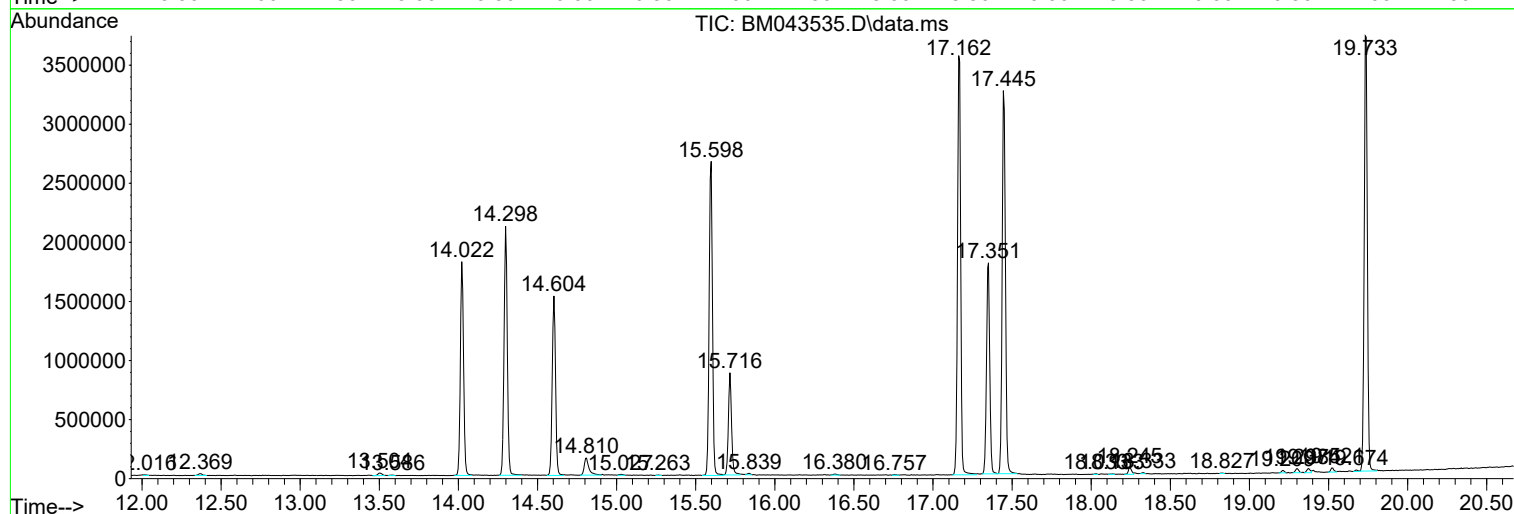
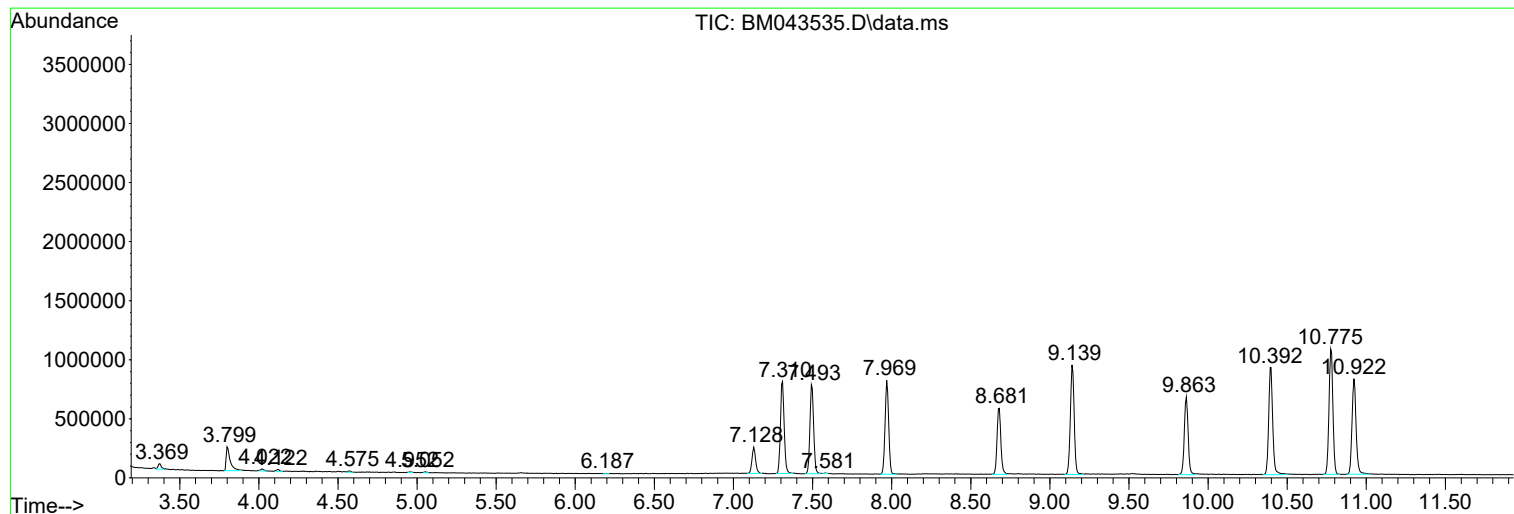
Sum of corrected areas: 54513058

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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



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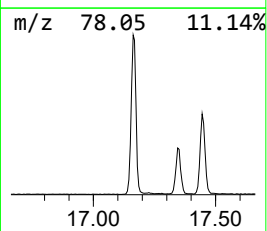
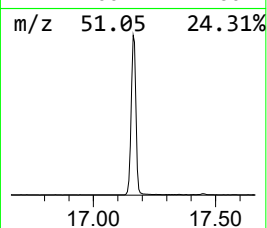
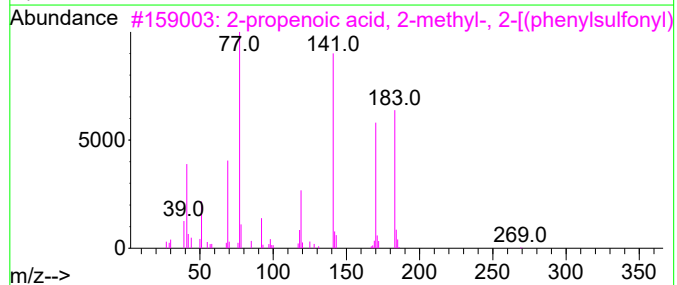
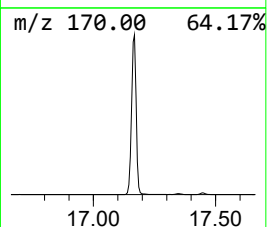
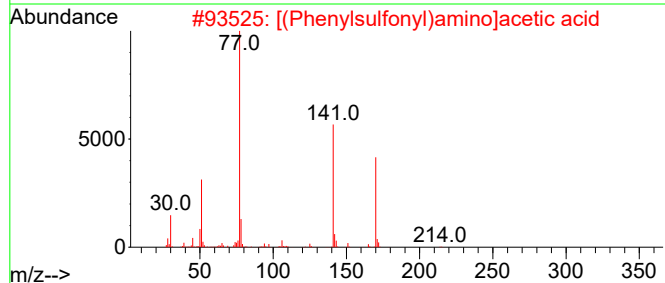
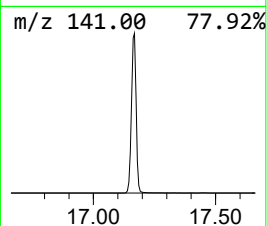
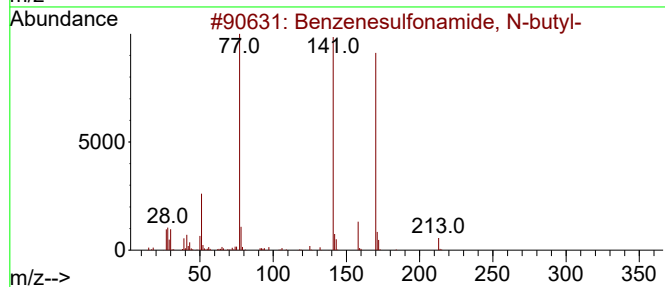
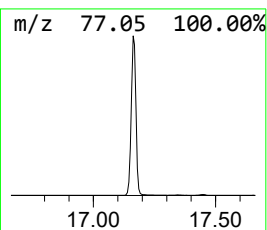
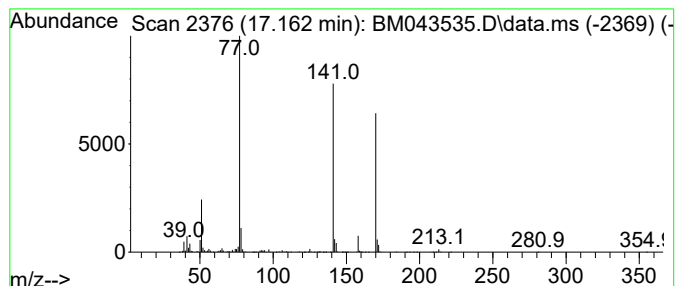
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Quant Title : SVOA CALIBRATION

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.
17.163	38.34 ng/ul	4890520	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	94
2			[(Phenylsulfonyl)amino]acetic acid	215	C8H9NO4S	005398-96-9	74
3			2-propenoic acid, 2-methyl-, 2-[...]	269	C12H15NO4S	1000401-71-8	74
4			N,N-Dichlorobenzenesulfonamide	225	C6H5Cl2NO2S	000473-29-0	64
5			N-Benzenesulfonyloxy-2,2-bis(tri...]	335	C10H7F6NO3S	055734-41-3	64



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzenesulfonam...	17.163	38.3	ng/ul	4890520	4	17.351	2551360	20.0