

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011725\
 Data File : BP023638.D
 Acq On : 17 Jan 2025 22:04
 Operator : RC/JU
 Sample : Q1086-15
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AF6

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-BP011425.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023638.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.287	29	34	40	rBV2	382766	656753	4.29%	0.510%
2	3.346	40	44	55	rVB	925102	1143760	7.47%	0.889%
3	3.587	81	85	92	rVB	29183	41933	0.27%	0.033%
4	3.717	103	107	118	rBV	413110	645087	4.22%	0.501%
5	4.052	159	164	169	rBV	29708	41292	0.27%	0.032%
6	4.146	173	180	187	rVB2	43274	72168	0.47%	0.056%
7	4.411	221	225	232	rVB9	10300	19728	0.13%	0.015%
8	4.505	237	241	245	rVV2	13064	20640	0.13%	0.016%
9	4.575	247	253	261	rVB6	19948	40283	0.26%	0.031%
10	4.952	312	317	323	rVB2	22558	36382	0.24%	0.028%
11	6.211	523	531	537	rVB	32233	54082	0.35%	0.042%
12	6.522	579	584	597	rBV6	12032	36067	0.24%	0.028%
13	6.964	652	659	678	rBV	517903	842192	5.50%	0.654%
14	7.140	681	689	698	rBV	1906380	3048952	19.92%	2.369%
15	7.340	715	723	738	rBV	2064188	3144616	20.55%	2.444%
16	7.805	794	802	810	rBV	1590643	2561740	16.74%	1.991%
17	7.928	816	823	830	rVB	229226	345292	2.26%	0.268%
18	8.487	910	918	931	rBV	1594210	2523682	16.49%	1.961%
19	8.946	988	996	1009	rBV	2084354	3371682	22.03%	2.620%
20	9.040	1009	1012	1018	rVB8	8075	15551	0.10%	0.012%
21	9.122	1021	1026	1030	rVV7	12163	19444	0.13%	0.015%
22	9.411	1070	1075	1082	rVB2	30146	51479	0.34%	0.040%
23	9.663	1104	1118	1130	rBV	1268841	2086935	13.64%	1.622%
24	10.199	1201	1209	1224	rBV	2660859	4203345	27.47%	3.266%
25	10.363	1232	1237	1248	rVB	59934	104234	0.68%	0.081%
26	10.581	1266	1274	1286	rVB	2116286	3644317	23.82%	2.832%
27	10.705	1287	1295	1311	rBV	2144424	3635752	23.76%	2.825%
28	11.399	1408	1413	1420	rVB3	32517	51260	0.33%	0.040%
29	11.899	1493	1498	1505	rBV5	10752	20885	0.14%	0.016%
30	12.163	1537	1543	1549	rBV	50319	77896	0.51%	0.061%
31	13.334	1736	1742	1746	rBV4	10715	16151	0.11%	0.013%
32	13.393	1748	1752	1757	rVB5	12627	19791	0.13%	0.015%
33	13.828	1818	1826	1833	rBV	4892747	6625167	43.30%	5.148%
34	13.975	1848	1851	1859	rVB10	10346	18537	0.12%	0.014%
35	14.116	1865	1875	1890	rBV	5683996	8506114	55.59%	6.610%
36	14.428	1918	1928	1935	rBV	3032147	4787337	31.29%	3.720%

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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-BP011425.MA.M
 Title : SVOA CALIBRATION

37	14.510	1938	1942	1946	rVB4	12580	21193	0.14%	0.016%
38	14.563	1946	1951	1952	rBV3	17842	22069	0.14%	0.017%
39	14.604	1952	1958	1979	rVB	469919	703019	4.59%	0.546%
40	14.846	1996	1999	2009	rVB	13085	21721	0.14%	0.017%
41	15.257	2064	2069	2074	rBV	31895	43474	0.28%	0.034%
42	15.434	2092	2099	2110	rBV	7571914	10490946	68.56%	8.152%
43	15.540	2111	2117	2136	rVV	1099839	1601560	10.47%	1.245%
44	15.681	2136	2141	2147	rVB2	33384	58012	0.38%	0.045%
45	15.804	2157	2162	2169	rVB	120977	173474	1.13%	0.135%
46	16.240	2231	2236	2244	rBV	59820	97458	0.64%	0.076%
47	16.381	2257	2260	2265	rVB2	12783	18946	0.12%	0.015%
48	17.022	2360	2369	2379	rBV	578894	866063	5.66%	0.673%
49	17.222	2396	2403	2413	rVB	4068889	5530630	36.14%	4.298%
50	17.322	2413	2420	2433	rBV	9282721	12298335	80.37%	9.557%
51	17.539	2452	2457	2460	rBV2	14647	24242	0.16%	0.019%
52	17.604	2464	2468	2471	rBV5	11252	16221	0.11%	0.013%
53	18.157	2559	2562	2570	rBV4	91718	138619	0.91%	0.108%
54	18.239	2573	2576	2582	rVB	12251	19458	0.13%	0.015%
55	18.481	2610	2617	2620	rBV9	16267	34453	0.23%	0.027%
56	18.634	2639	2643	2649	rVB9	16007	29037	0.19%	0.023%
57	18.875	2681	2684	2688	rVB5	13916	16558	0.11%	0.013%
58	19.163	2728	2733	2741	rBV	217205	297894	1.95%	0.231%
59	19.245	2742	2747	2752	rBV	92908	131308	0.86%	0.102%
60	19.310	2754	2758	2764	rVV	86808	118328	0.77%	0.092%
61	19.492	2785	2789	2795	rBV2	68489	97519	0.64%	0.076%
62	19.692	2817	2823	2835	rBV	11347078	14716672	96.17%	11.436%
63	21.680	3155	3161	3175	rBV2	3833877	6223114	40.67%	4.836%
64	21.863	3188	3192	3202	rVB	269178	431795	2.82%	0.336%
65	24.880	3694	3705	3724	rBV	5869357	15302304	100.00%	11.891%
66	25.110	3734	3744	3759	rVB	2564873	6615323	43.23%	5.140%

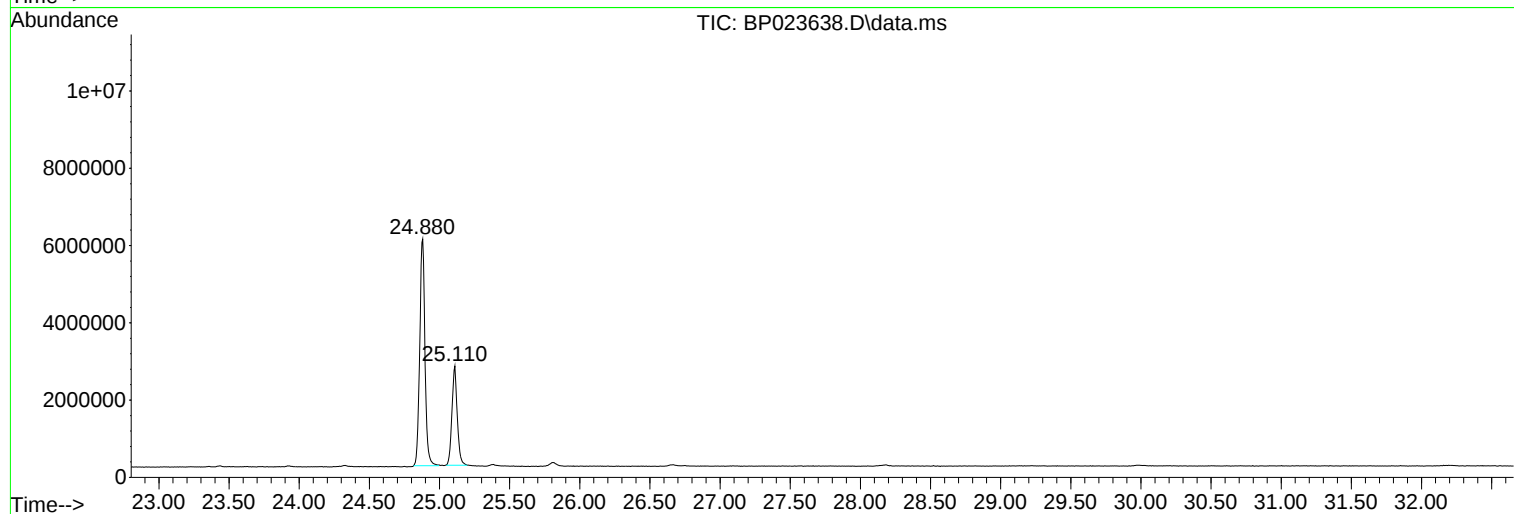
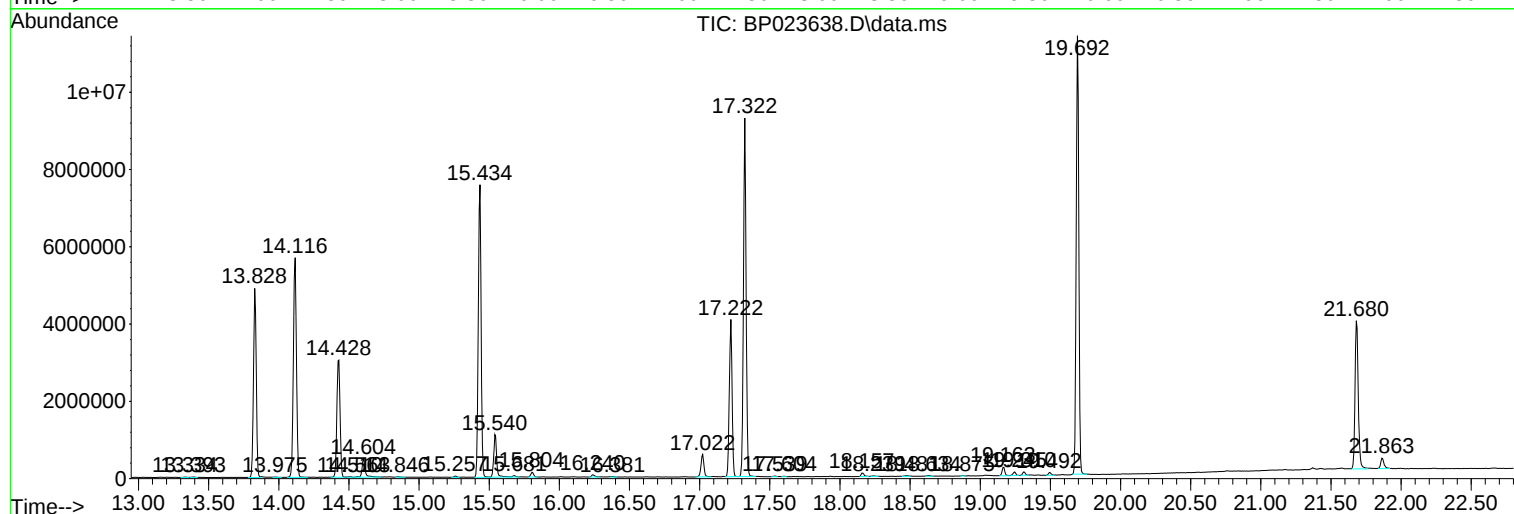
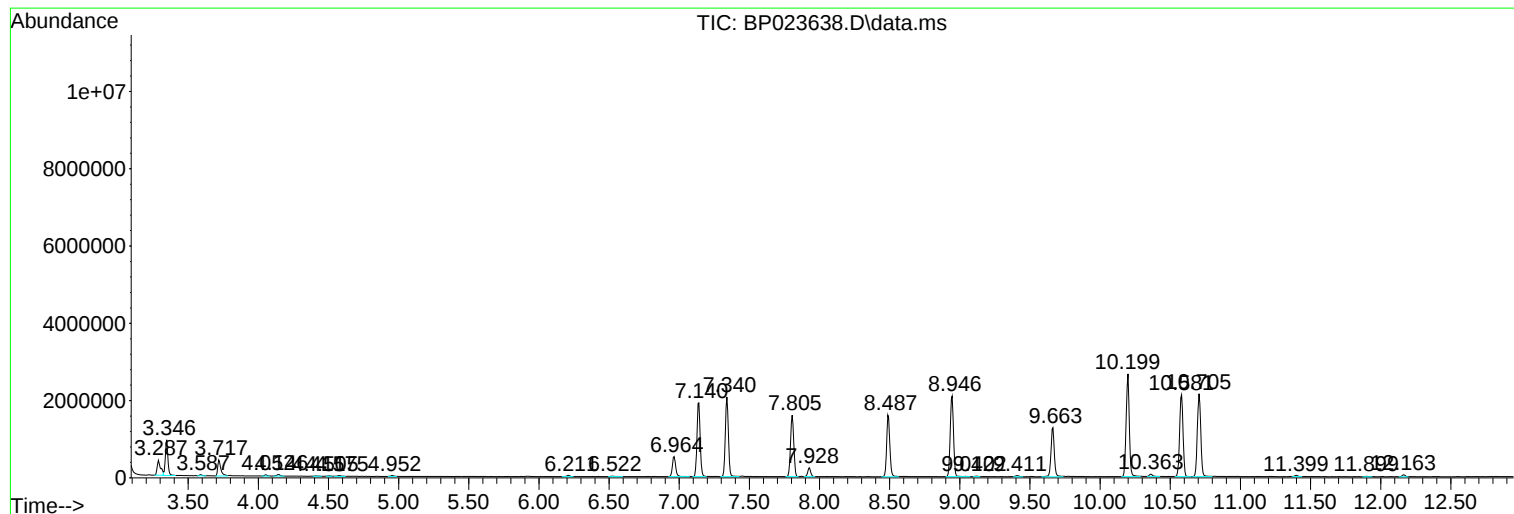
Sum of corrected areas: 128690271

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
Quant Title : SVOA CALIBRATION

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



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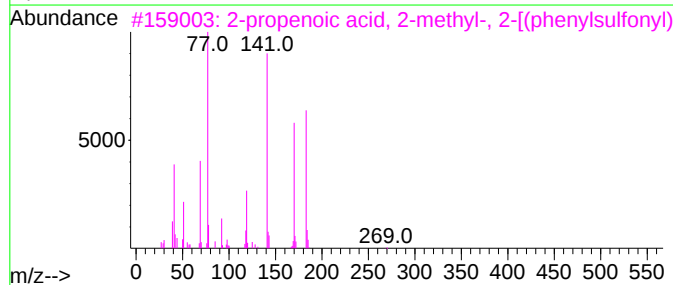
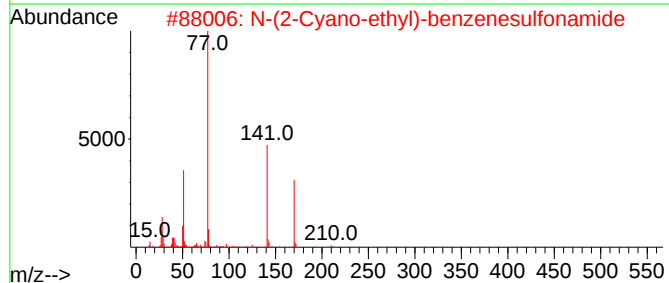
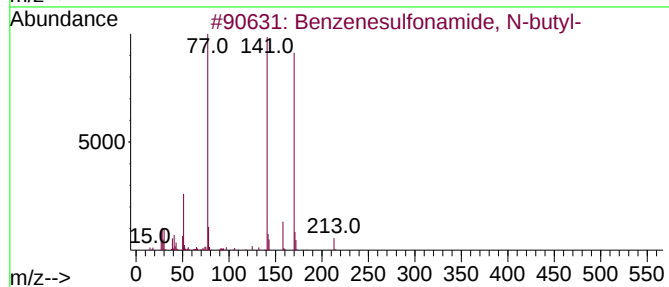
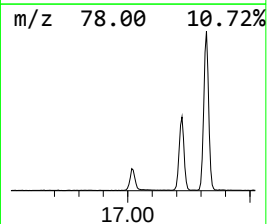
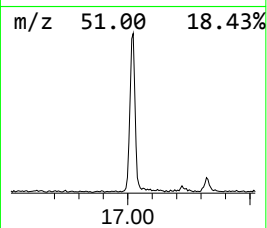
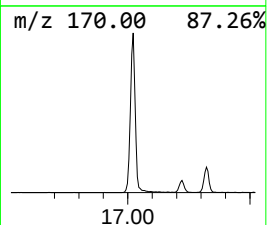
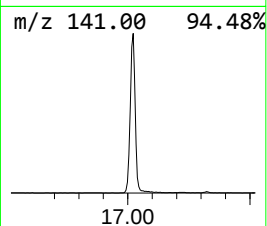
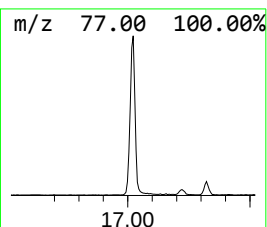
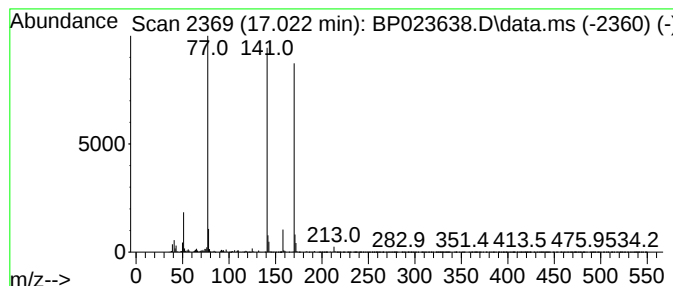
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzenesulfonamide, N-butyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.022	3.13 ng/ul	866063	Phenanthrene-d10	17.222

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	91
2			N-(2-Cyano-ethyl)-benzenesulfona...	210	C9H10N2O2S	002619-21-8	74
3			2-propenoic acid, 2-methyl-, 2-[...]	269	C12H15NO4S	1000401-71-8	64
4			N-Octylbenzenesulfonamide	269	C14H23NO2S	016358-32-0	50
5			N-Benzenesulfonyloxy-2,2-bis(tri...	335	C10H7F6NO3S	055734-41-3	50



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzenesulfonam...	17.022	3.1	ng/u1	866063	4	17.222	5530630	20.0