

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090823\
 Data File : BG058941.D
 Acq On : 9 Sep 2023 8:03
 Operator : MA/JU
 Sample : 04230-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BH490

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_G\Methods\SFAM-EPA-BG090823.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG058941.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.610	16	21	23	rBV3	19392	30446	0.19%	0.077%
2	4.075	96	100	110	rBV	58496	120855	0.77%	0.307%
3	7.435	667	672	681	rVB2	66481	111096	0.71%	0.283%
4	7.647	701	708	718	rVB	206249	367332	2.34%	0.934%
5	7.841	735	741	748	rVB	179020	322229	2.05%	0.820%
6	8.328	818	824	832	rVB	149494	261337	1.66%	0.665%
7	9.004	932	939	946	rVB2	177601	323314	2.06%	0.823%
8	9.521	1021	1027	1038	rVB	261460	479664	3.06%	1.220%
9	10.244	1141	1150	1158	rBV2	217091	393354	2.51%	1.001%
10	10.767	1232	1239	1246	rVB	373416	647447	4.12%	1.647%
11	11.172	1301	1308	1316	rVB2	232677	418702	2.67%	1.065%
12	11.313	1325	1332	1340	rVB2	361943	677552	4.32%	1.724%
13	12.723	1569	1572	1578	rVB3	9816	17354	0.11%	0.044%
14	13.793	1750	1754	1757	rVB3	12653	16897	0.11%	0.043%
15	14.351	1842	1849	1857	rBV	1024767	1459874	9.30%	3.714%
16	14.656	1893	1901	1908	rBV2	961347	1537460	9.79%	3.911%
17	14.950	1942	1951	1961	rVB	490234	804892	5.13%	2.048%
18	15.115	1971	1979	1985	rBV2	89510	145337	0.93%	0.370%
19	15.943	2111	2120	2129	rBV	1384865	2138867	13.63%	5.441%
20	16.043	2129	2137	2142	rBV	384610	557714	3.55%	1.419%
21	16.178	2155	2160	2169	rVB2	11483	18315	0.12%	0.047%
22	17.541	2375	2392	2396	rBV	6766988	15696921	100.00%	39.933%
23	17.706	2413	2420	2427	rVB	797293	1212725	7.73%	3.085%
24	17.806	2430	2437	2444	rVB2	1744027	2627043	16.74%	6.683%
25	18.528	2554	2560	2565	rBV3	52819	76891	0.49%	0.196%
26	19.515	2723	2728	2731	rBV3	22669	33528	0.21%	0.085%
27	19.598	2735	2742	2745	rBV2	26437	38468	0.25%	0.098%
28	19.639	2745	2749	2753	rVV2	29406	37792	0.24%	0.096%
29	19.703	2757	2760	2767	rVB2	34566	46527	0.30%	0.118%
30	19.792	2769	2775	2781	rBV6	32647	50996	0.32%	0.130%
31	19.944	2792	2801	2804	rBV7	18546	29752	0.19%	0.076%
32	20.074	2817	2823	2832	rVB	2176688	3139105	20.00%	7.986%
33	20.990	2974	2979	2982	rBV6	12487	18253	0.12%	0.046%
34	21.589	3078	3081	3085	rBV5	16889	26853	0.17%	0.068%
35	22.030	3147	3156	3167	rVV2	650610	1158622	7.38%	2.948%
36	22.124	3167	3172	3178	rVV2	42843	74741	0.48%	0.190%

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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_G\Methods\SFAM-EPA-BG090823.MA.M
Title : SVOA CALIBRATION

37	22.177	3178	3181	3185	rVB6	11984	17134	0.11%	0.044%
38	25.350	3709	3721	3735	rBV2	892871	2763594	17.61%	7.031%
39	25.602	3753	3764	3776	rVB	407135	1244031	7.93%	3.165%
40	26.149	3849	3857	3864	rBV5	36800	105296	0.67%	0.268%
41	26.883	3974	3982	3983	rBV5	21514	44108	0.28%	0.112%
42	28.440	4245	4247	4250	rVB3	12940	15979	0.10%	0.041%

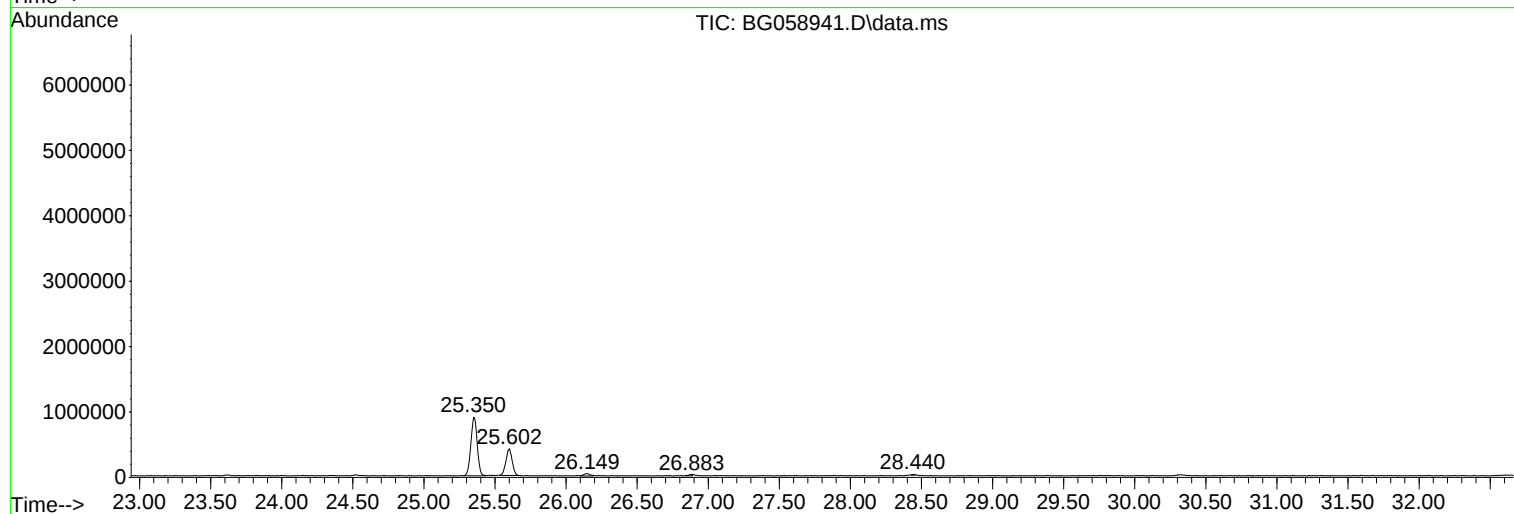
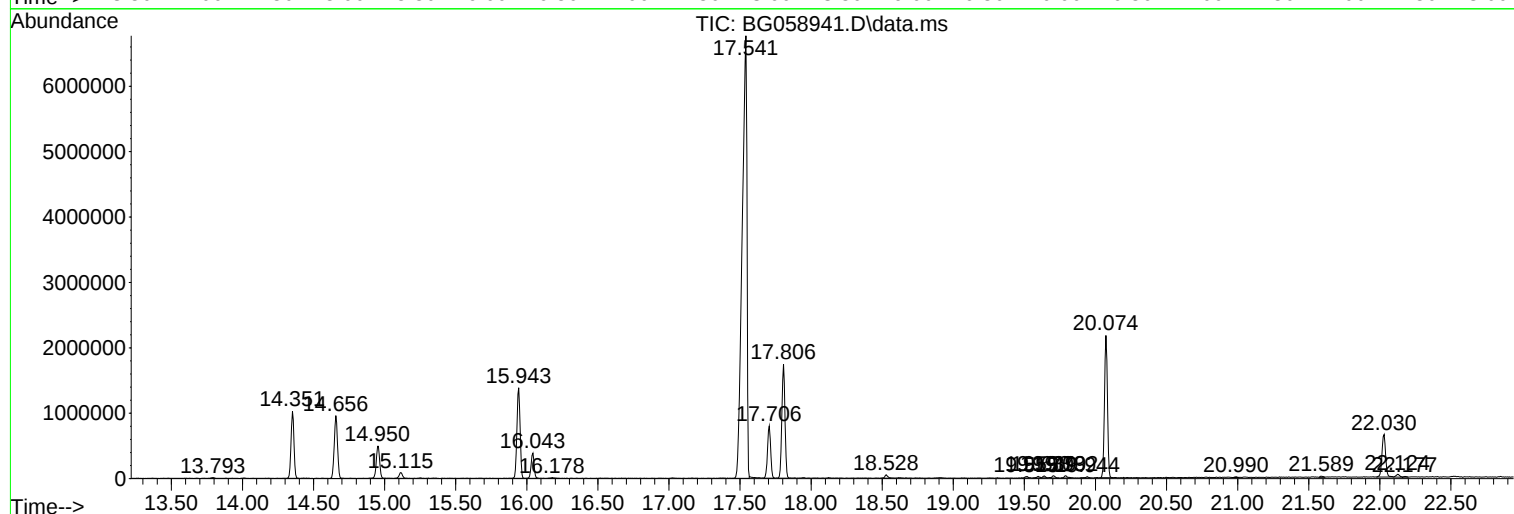
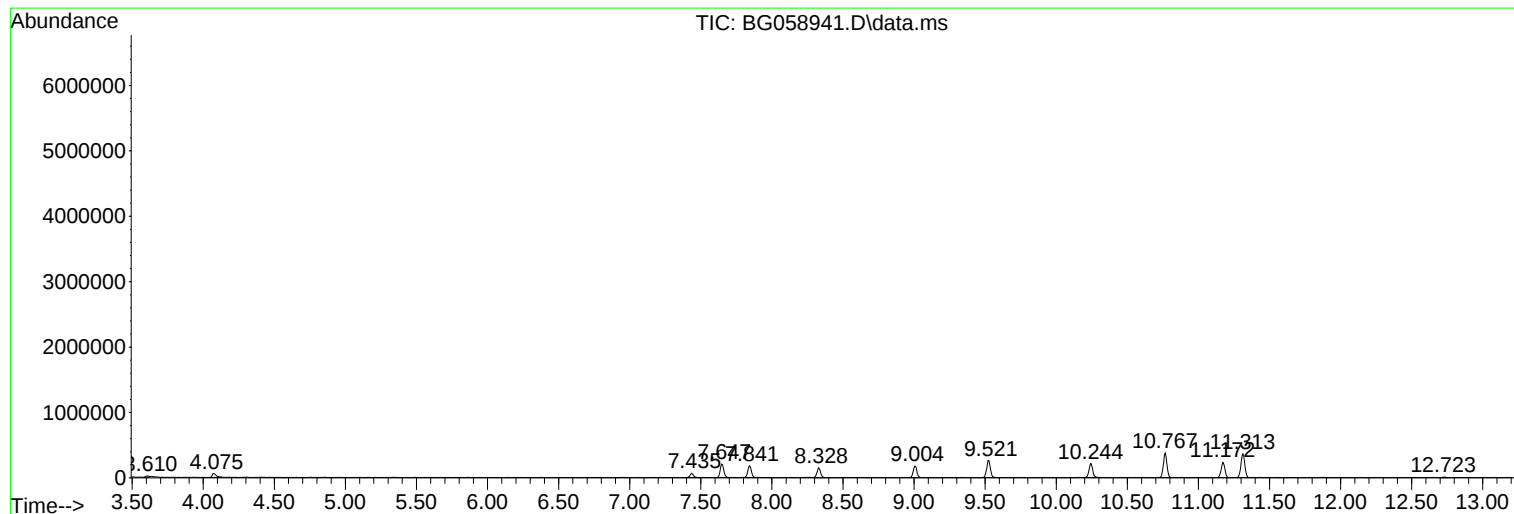
Sum of corrected areas: 39308397

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090823.MA.M
Quant Title : SVOA CALIBRATION

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



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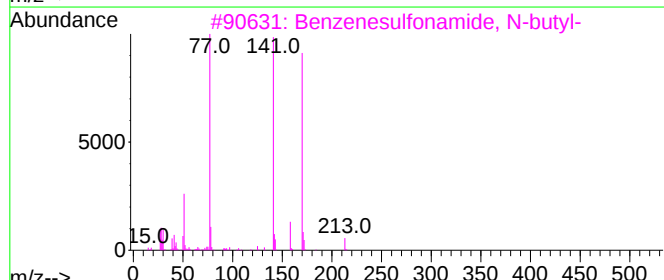
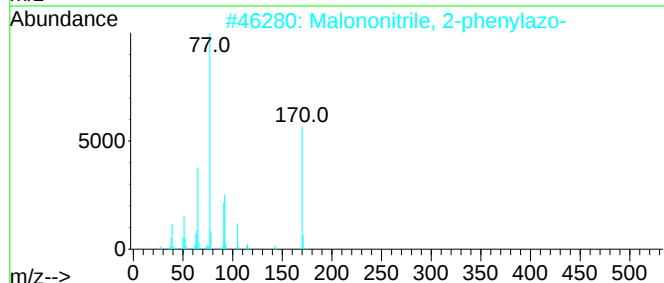
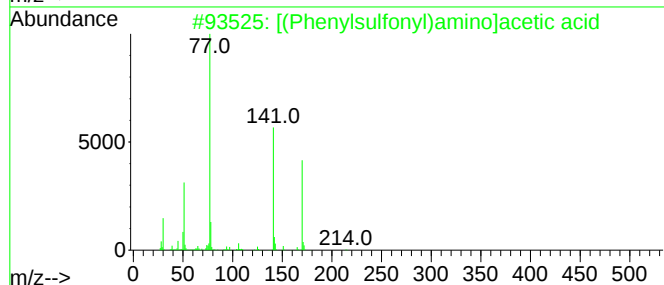
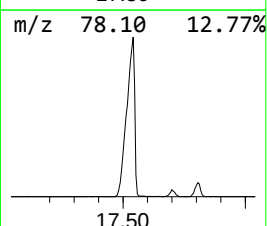
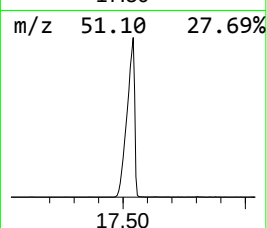
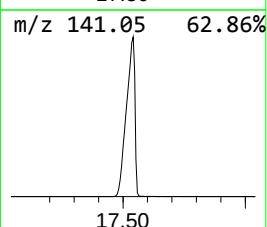
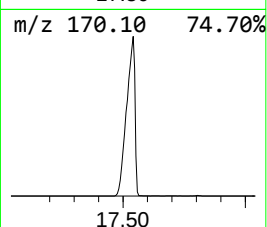
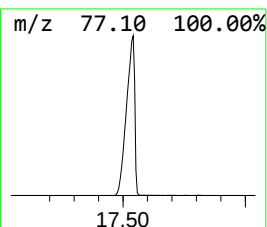
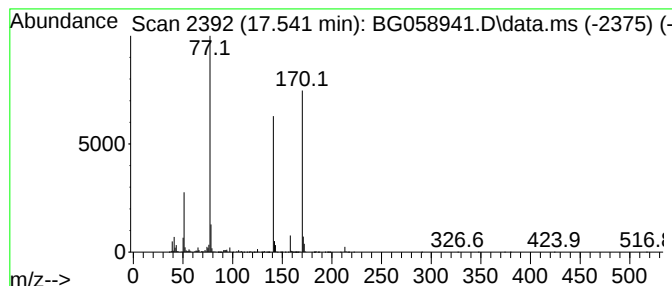
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 [(Phenylsulfonyl)amino]acet... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.541	258.87 ng/ul	15696900	Phenanthrene-d10	17.706

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		[(Phenylsulfonyl)amino]acetic acid	215	C8H9NO4S	005398-96-9	53
2		Malononitrile, 2-phenylazo-	170	C9H6N4	006017-21-6	53
3		Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	49
4		(Phenylsulfonyl)acetonitrile	181	C8H7NO2S	007605-28-9	47
5		1-[Phenylsulfonyl]-1-nitropropane	229	C9H11NO4S	021272-84-4	47



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
[(Phenylsulfony...	17.541	258.9	ng/ul	15696900	4	17.706	1212730	20.0