

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
 Data File : BP019182.D
 Acq On : 30 Dec 2023 00:56
 Operator : MA/JU
 Sample : 06021-04
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH7Y6

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-BP122723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019182.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.199	15	19	23	rBV	194808	256989	1.23%	0.346%
2	3.234	23	25	30	rVB	55672	64764	0.31%	0.087%
3	3.658	93	97	110	rBV	205912	310896	1.49%	0.419%
4	3.975	148	151	158	rVB	20705	26838	0.13%	0.036%
5	4.422	223	227	232	rVB3	24485	36268	0.17%	0.049%
6	6.975	655	661	671	rVB	207344	326874	1.56%	0.441%
7	7.140	682	689	698	rBV	704384	1088027	5.21%	1.467%
8	7.328	714	721	736	rBV	671888	1078179	5.16%	1.453%
9	7.799	793	801	809	rBV	708063	1105457	5.29%	1.490%
10	8.516	915	923	936	rBV	562191	916232	4.38%	1.235%
11	8.963	991	999	1008	rBV	780645	1250341	5.98%	1.685%
12	9.681	1114	1121	1134	rBV	569685	948080	4.54%	1.278%
13	10.222	1204	1213	1230	rBV	906332	1578075	7.55%	2.127%
14	10.593	1268	1276	1284	rBV	894981	1491211	7.13%	2.010%
15	10.740	1292	1301	1320	rBV	701821	1277601	6.11%	1.722%
16	12.193	1541	1548	1556	rBV2	16435	31493	0.15%	0.042%
17	13.345	1738	1744	1750	rVV2	23979	39788	0.19%	0.054%
18	13.863	1821	1832	1840	rBV	1872157	2573088	12.31%	3.468%
19	14.134	1866	1878	1887	rBV	2030195	2972724	14.22%	4.007%
20	14.445	1922	1931	1944	rBV	1357139	2107704	10.08%	2.841%
21	14.663	1961	1968	1984	rBV	124115	268755	1.29%	0.362%
22	15.439	2092	2100	2108	rBV	2892156	3986655	19.07%	5.374%
23	15.563	2115	2121	2136	rBV	744454	1049606	5.02%	1.415%
24	16.222	2227	2233	2241	rBV6	15140	31129	0.15%	0.042%
25	17.039	2360	2372	2377	rBV	10520954	20900270	100.00%	28.173%
26	17.157	2388	2392	2394	rBV3	43314	56344	0.27%	0.076%
27	17.198	2394	2399	2407	rVB	1827238	2498372	11.95%	3.368%
28	17.298	2409	2416	2427	rBV	3345370	4641589	22.21%	6.257%
29	18.092	2547	2551	2560	rBV2	82674	124734	0.60%	0.168%
30	18.169	2560	2564	2569	rVB2	16555	24875	0.12%	0.034%
31	19.028	2706	2710	2716	rBV	86412	107513	0.51%	0.145%
32	19.116	2719	2725	2732	rBV3	52017	89135	0.43%	0.120%
33	19.192	2733	2738	2740	rBV	46806	68221	0.33%	0.092%
34	19.216	2740	2742	2750	rVB6	32606	50080	0.24%	0.068%
35	19.333	2758	2762	2767	rBV3	60198	80209	0.38%	0.108%
36	19.545	2792	2798	2810	rVV	4409640	5922996	28.34%	7.984%

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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

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37	21.298	3091	3096	3104	rBV	2478417	3215379	15.38%	4.334%
38	21.398	3109	3113	3118	rVV	260522	318134	1.52%	0.429%
39	23.510	3465	3472	3488	rVB2	3563836	7115716	34.05%	9.592%
40	23.663	3492	3498	3510	rVB	1868827	3570276	17.08%	4.813%
41	23.927	3539	3543	3551	rVB	321163	585912	2.80%	0.790%

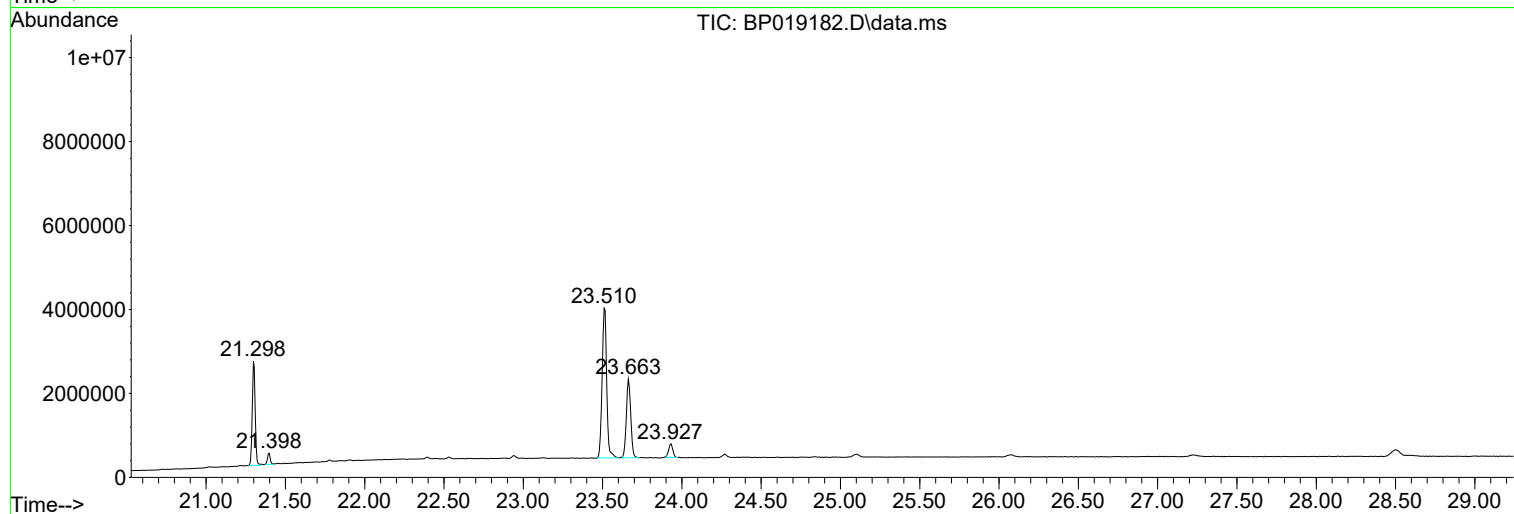
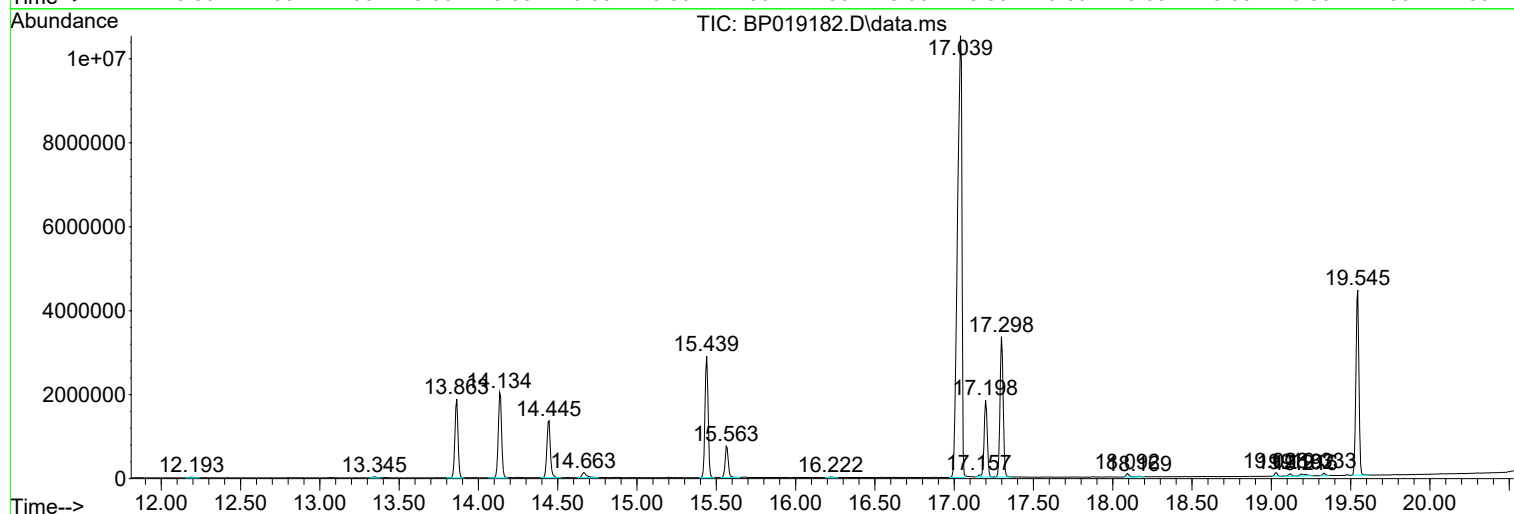
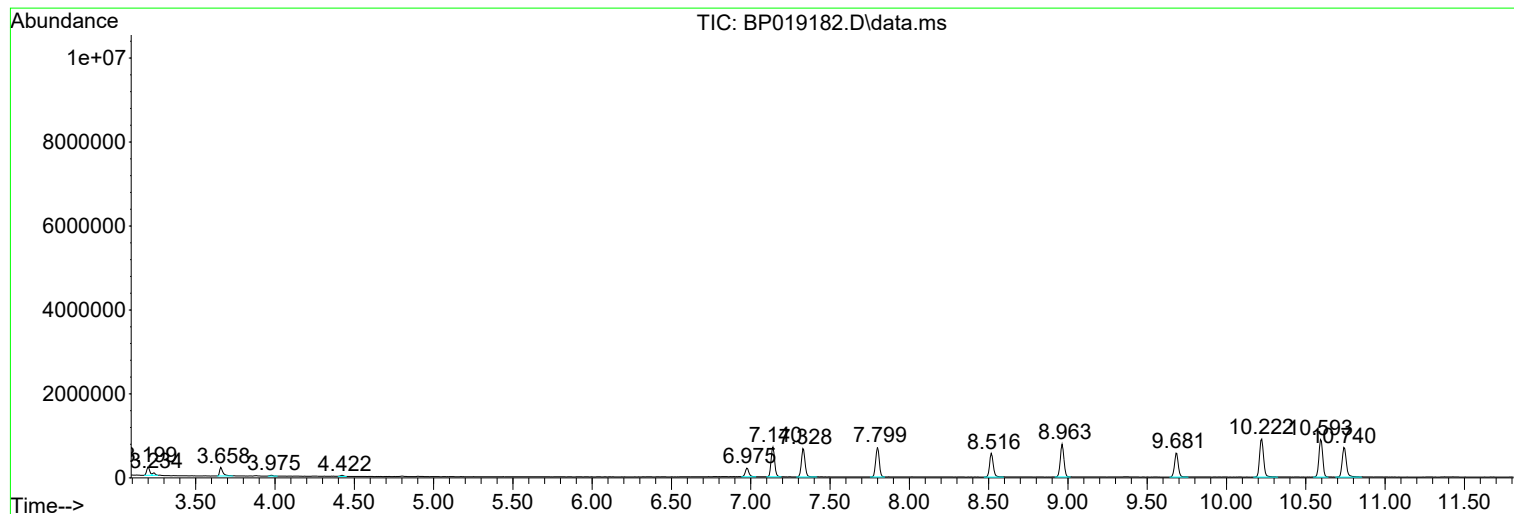
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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



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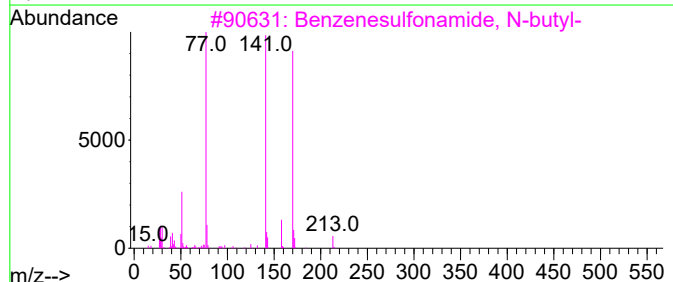
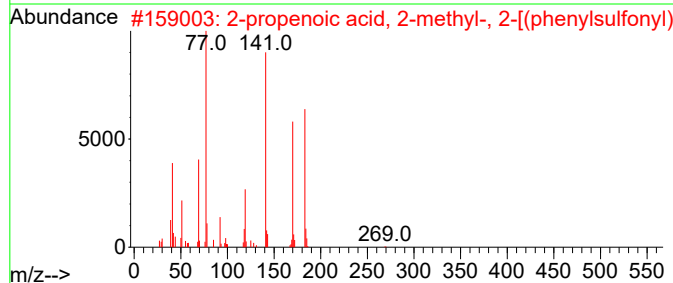
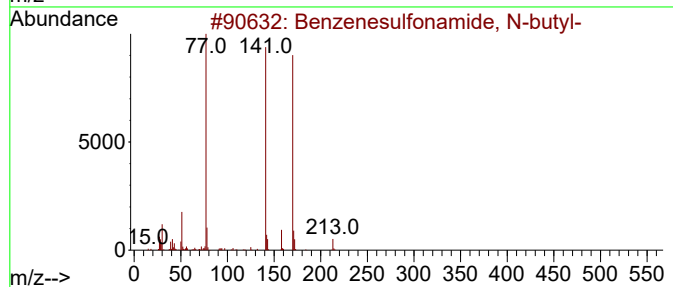
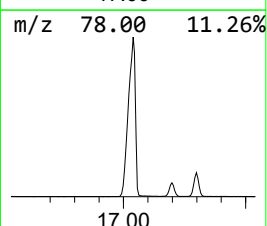
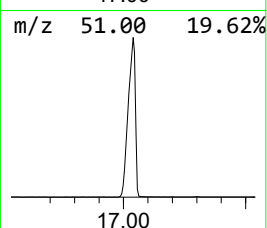
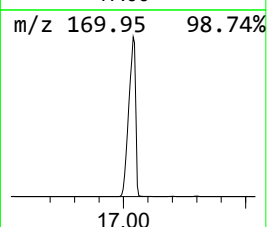
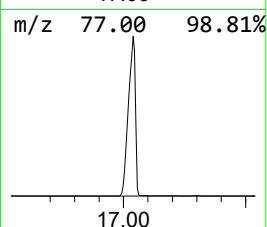
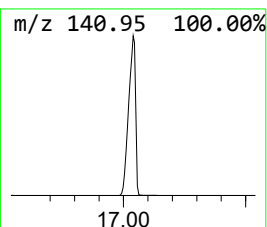
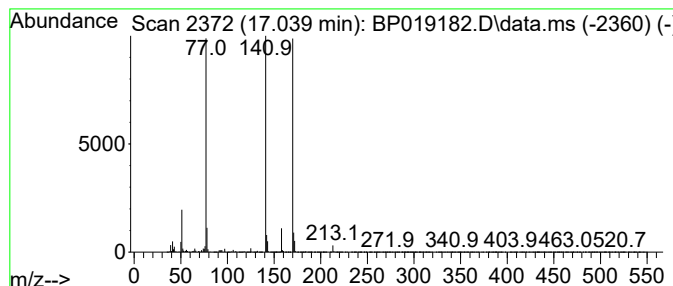
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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.
17.039	167.31 ng/ul	20900300	Phenanthrene-d10	17.198

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	90
2			2-propenoic acid, 2-methyl-, 2-[...]	269	C12H15NO4S	1000401-71-8	64
3			Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	53
4			N-Benzenesulfonyloxy-2,2-bis(tri...]	335	C10H7F6N03S	055734-41-3	50
5			N-(2-Pyrazinyl)benzenesulfonamide	235	C10H9N3O2S	007471-20-7	47



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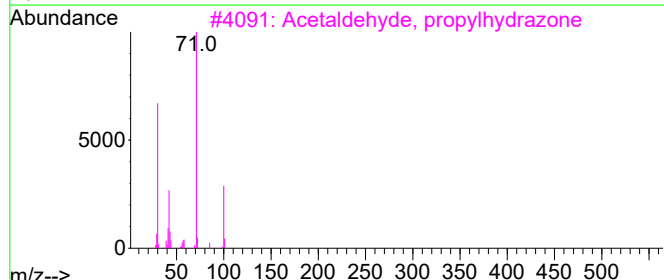
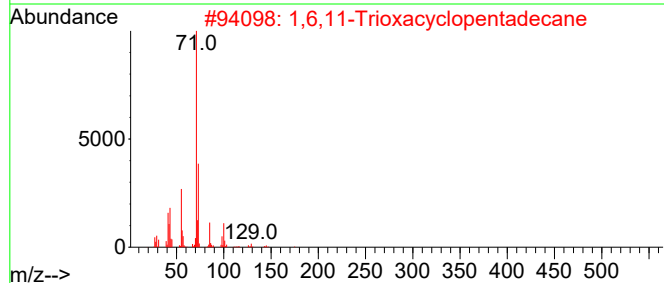
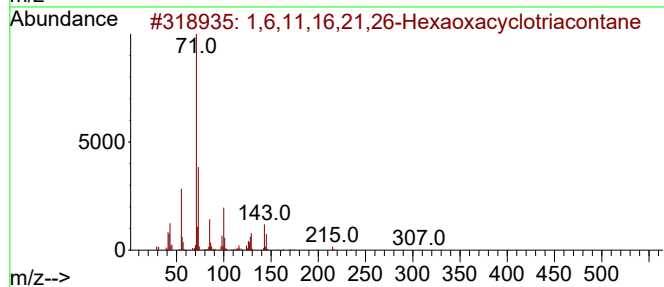
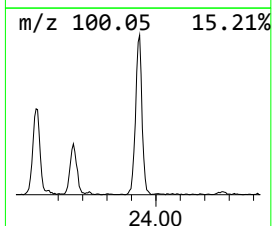
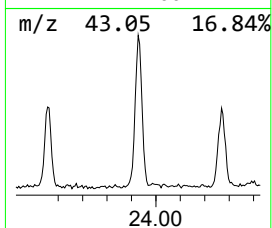
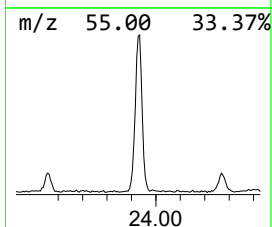
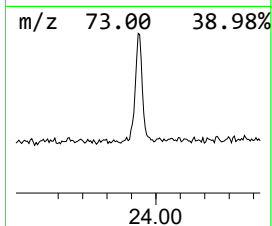
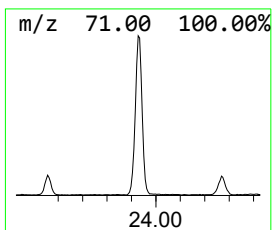
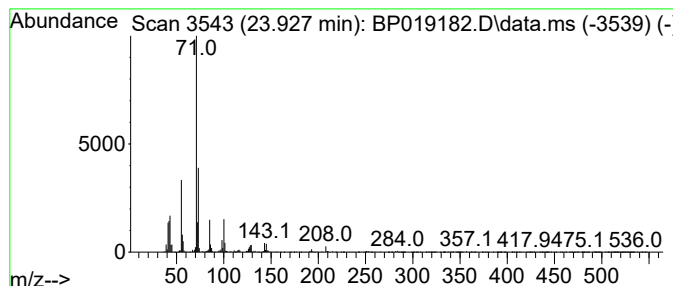
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TIC Library : C:\DATABASE\NIST20.L
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Peak Number 2 1,6,11,16,21,26-Hexaoxacycl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.927	3.28 ng/ul	585912	Perylene-d12	23.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,6,11,16,21,26-Hexaoxacyclotria...	432	C24H48O6	064001-05-4	90		
2	1,6,11-Trioxacyclopentadecane	216	C12H24O3	000295-63-6	83		
3	Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	58		
4	Acrolein,dimethyl acetal	102	C5H10O2	006044-68-4	50		
5	1,6,11,16,21,26,31-Heptaoxacyclo...	504	C28H56O7	066055-34-3	50		



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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...	17.039	167.3	ng/ul	20900300	4	17.198	2498370	20.0
1,6,11,16,21,26...	23.927	3.3	ng/ul	585912	6	23.663	3570280	20.0