

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
 Data File : BN033394.D
 Acq On : 15 Aug 2024 11:02
 Operator : MA/JU
 Sample : PB162649BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK649

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_N\Methods\SFAM-EPA-BN081424.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN033394.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.158	25	29	42	rVB	85134	127085	1.48%	0.132%
2	3.540	90	94	116	rBV	1058415	1751699	20.33%	1.816%
3	3.864	145	149	153	rBV	19754	26098	0.30%	0.027%
4	6.569	606	609	617	rBV9	3978	8887	0.10%	0.009%
5	6.746	632	639	651	rBV	1491429	2352835	27.31%	2.439%
6	6.916	660	668	678	rVB	1246417	1870410	21.71%	1.939%
7	7.105	693	700	712	rBV	1500732	2278983	26.45%	2.362%
8	7.569	772	779	787	rBV	1432220	2245441	26.06%	2.327%
9	8.269	889	898	908	rBV	1927028	2999425	34.82%	3.109%
10	8.710	963	973	984	rBV	1437184	2342254	27.19%	2.428%
11	9.422	1083	1094	1105	rBV	1113110	1803537	20.93%	1.869%
12	9.957	1177	1185	1195	rBV	1893355	3014537	34.99%	3.124%
13	10.334	1241	1249	1258	rBV	2161860	3554448	41.26%	3.684%
14	10.469	1264	1272	1282	rBV	1804518	2976517	34.55%	3.085%
15	11.928	1515	1520	1526	rVB2	36628	58381	0.68%	0.061%
16	13.628	1802	1809	1817	rBV	3206031	4439714	51.53%	4.601%
17	13.898	1847	1855	1862	rBV	3747551	5462274	63.40%	5.661%
18	14.210	1901	1908	1915	rBV	3285912	4891359	56.78%	5.070%
19	14.281	1918	1920	1926	rVB7	7755	11795	0.14%	0.012%
20	14.410	1934	1942	1953	rBV	1769792	2436825	28.29%	2.526%
21	14.563	1964	1968	1974	rVV6	5873	9067	0.11%	0.009%
22	14.628	1977	1979	1985	rVB5	8542	11855	0.14%	0.012%
23	15.204	2070	2077	2085	rBV	4695848	6681946	77.56%	6.925%
24	15.322	2091	2097	2109	rBV	1340939	1755917	20.38%	1.820%
25	15.445	2113	2118	2123	rVB	20918	31935	0.37%	0.033%
26	15.722	2159	2165	2170	rBV8	4863	8948	0.10%	0.009%
27	15.992	2206	2211	2216	rBV4	15215	23800	0.28%	0.025%
28	16.745	2335	2339	2342	rBV2	7842	12307	0.14%	0.013%
29	16.957	2368	2375	2384	rBV	4150239	5834457	67.72%	6.047%
30	17.057	2385	2392	2404	rBV2	5369428	7410420	86.02%	7.680%
31	17.339	2436	2440	2449	rBV2	24280	46389	0.54%	0.048%
32	17.645	2490	2492	2500	rVB9	7969	15792	0.18%	0.016%
33	17.886	2529	2533	2538	rBV6	19385	29839	0.35%	0.031%
34	18.175	2577	2582	2588	rBV9	6642	14409	0.17%	0.015%
35	18.922	2705	2709	2716	rVB	70446	95066	1.10%	0.099%
36	18.992	2716	2721	2728	rBV	65674	98167	1.14%	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_N\Methods\SFAM-EPA-BN081424.MA.M
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37	19.180	2750	2753	2756	rBV4	13958	16125	0.19%	0.017%
38	19.357	2777	2783	2789	rBV	6374041	8615005	100.00%	8.929%
39	19.422	2789	2794	2808	rVB	442282	709662	8.24%	0.736%
40	20.398	2957	2960	2963	rVB5	18751	20226	0.23%	0.021%
41	21.192	3089	3095	3107	rBV2	4693805	6399775	74.29%	6.633%
42	23.833	3534	3544	3557	rBV2	3418746	7995367	92.81%	8.287%
43	24.021	3567	3576	3589	rVB	2556134	5995975	69.60%	6.214%

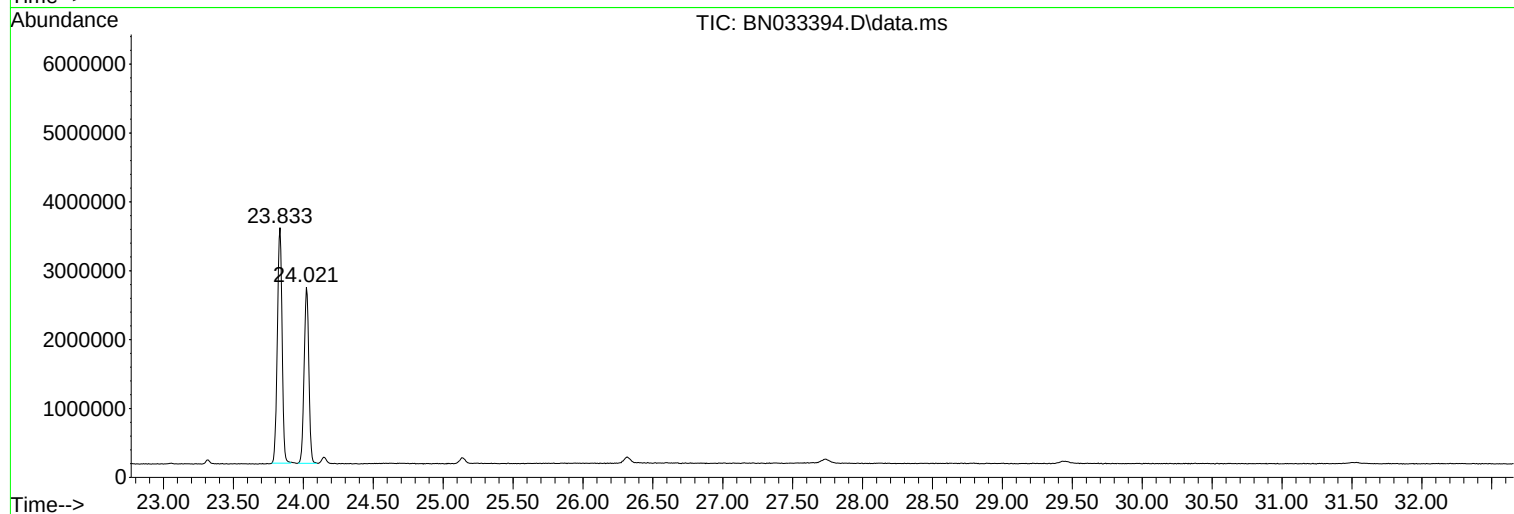
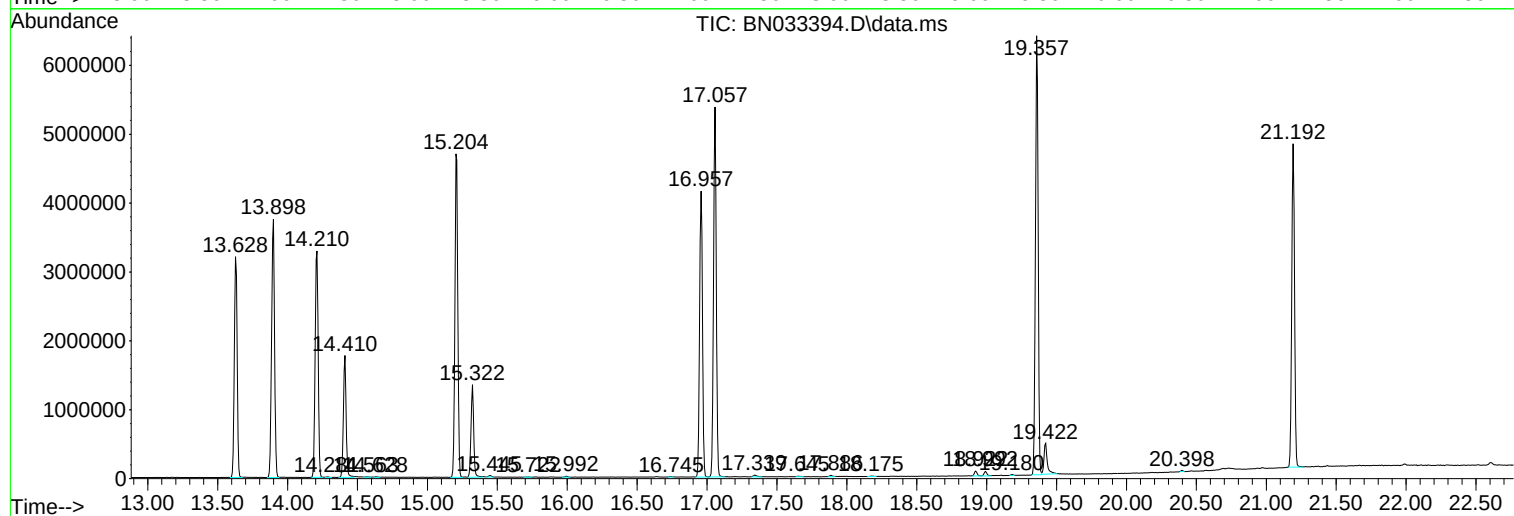
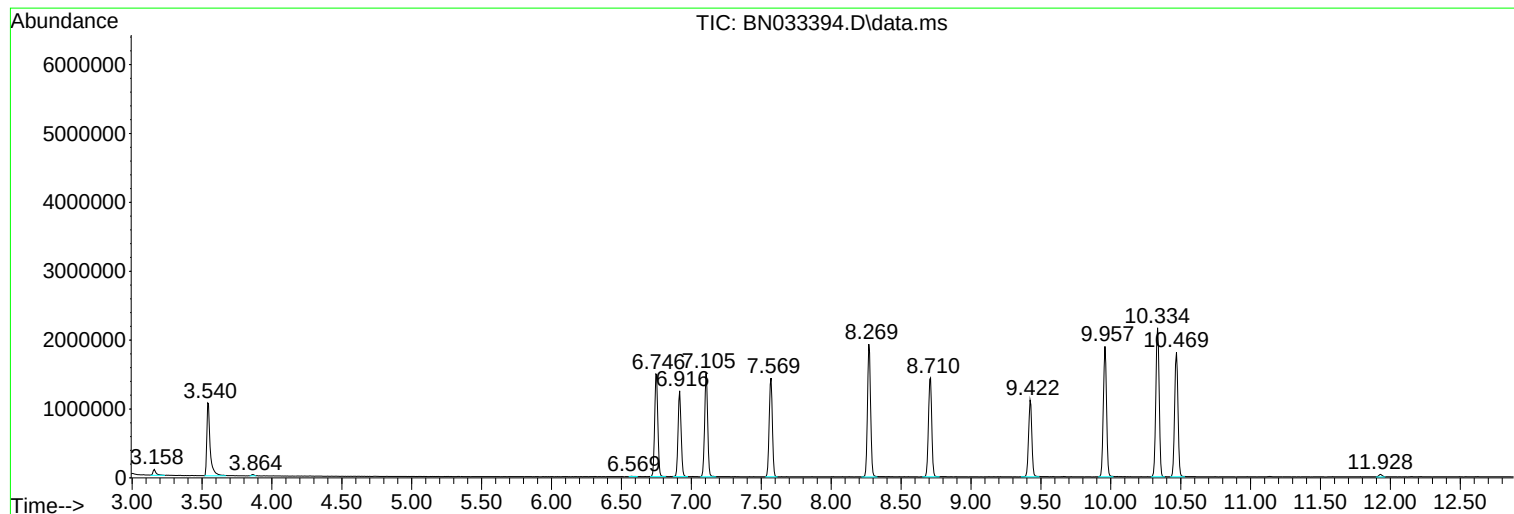
Sum of corrected areas: 96484953

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Quant Title : SVOA CALIBRATION

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



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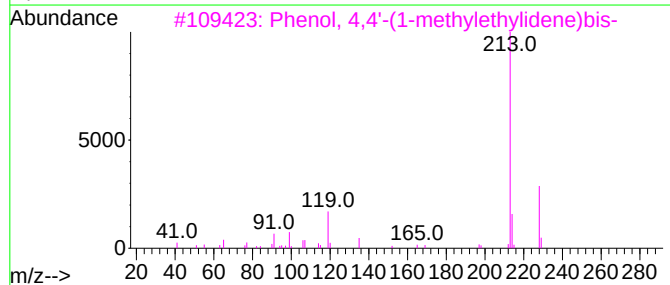
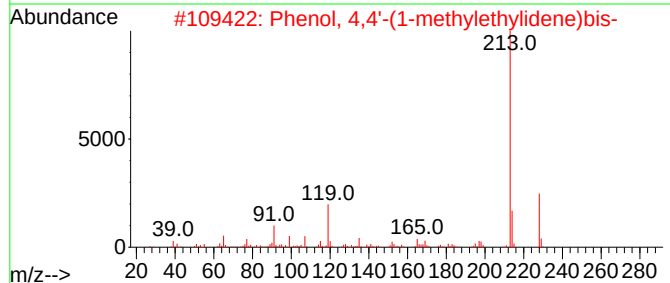
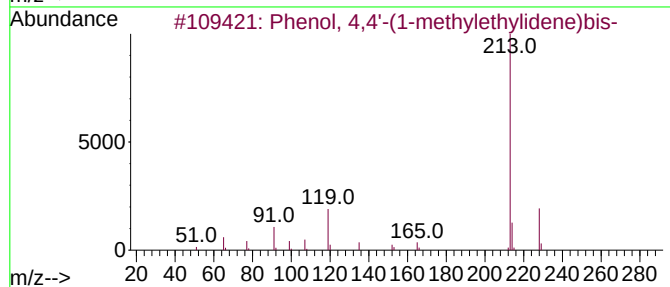
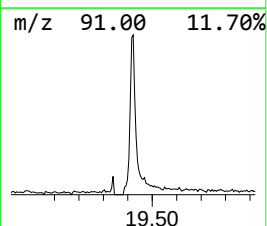
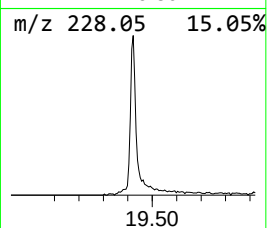
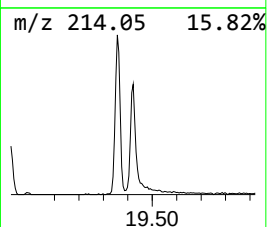
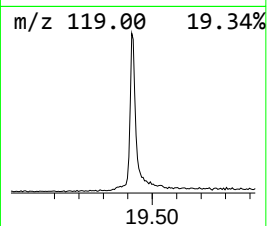
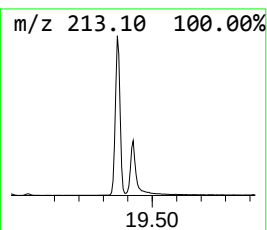
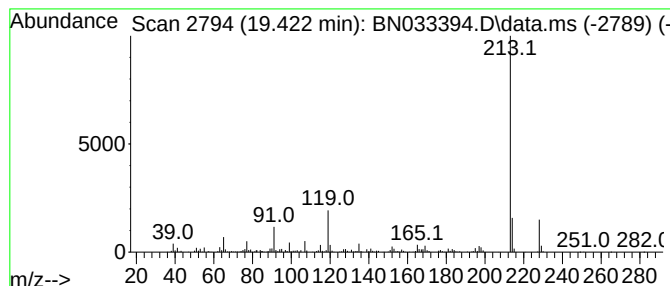
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenol, 4,4'-(1-methylethyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.422	2.22 ng/ul	709662	Chrysene-d12	21.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	96
2			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	93
3			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	90
4			4,4'-(1,3-Dimethylbutylidene)bis...	270	C18H22O2	006807-17-6	74
5			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	72



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 4,4'-(1...	19.422	2.2	ng/u1	709662	5	21.192	6399780	20.0