

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091625\
 Data File : BG064402.D
 Acq On : 17 Sep 2025 2:58
 Operator : RC/JU
 Sample : Q3084-18
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PMW-2(20250911)

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-BG091125.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG064402.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	28	34	41	rBV	234266	422369	38.17%	3.647%
2	3.352	41	45	55	rVB	54891	96763	8.75%	0.835%
3	3.722	103	108	116	rBV3	35163	59517	5.38%	0.514%
4	3.980	149	152	154	rBV3	1275	1514	0.14%	0.013%
5	4.057	160	165	170	rVB5	2873	4453	0.40%	0.038%
6	4.962	315	319	323	rBV3	1446	2919	0.26%	0.025%
7	5.173	351	355	359	rBV2	10240	15289	1.38%	0.132%
8	5.960	486	489	493	rVB	1735	1691	0.15%	0.015%
9	6.201	528	530	534	rVB2	2012	2389	0.22%	0.021%
10	6.554	586	590	596	rBV2	2884	4348	0.39%	0.038%
11	7.012	662	668	675	rBV	33998	54876	4.96%	0.474%
12	7.171	689	695	702	rVB	69616	116056	10.49%	1.002%
13	7.376	724	730	738	rVV	98769	166056	15.01%	1.434%
14	7.453	738	743	749	rVV	9449	14902	1.35%	0.129%
15	7.729	787	790	795	rBV2	6950	11497	1.04%	0.099%
16	7.840	803	809	814	rBV	78807	142938	12.92%	1.234%
17	7.958	826	829	835	rVB2	2353	3937	0.36%	0.034%
18	8.546	922	929	936	rBV	93028	155453	14.05%	1.342%
19	8.822	970	976	978	rBV2	2134	2950	0.27%	0.025%
20	8.992	998	1005	1014	rVB	120067	208474	18.84%	1.800%
21	9.562	1097	1102	1105	rVB	1276	1840	0.17%	0.016%
22	9.662	1116	1119	1120	rBV	1484	1589	0.14%	0.014%
23	9.715	1121	1128	1135	rVV	110178	193091	17.45%	1.667%
24	10.255	1212	1220	1227	rBV	148969	268316	24.25%	2.317%
25	10.496	1256	1261	1267	rVB4	9315	14866	1.34%	0.128%
26	10.631	1276	1284	1290	rBV	108060	187836	16.98%	1.622%
27	10.761	1299	1306	1319	rVB	139112	248145	22.43%	2.143%
28	11.730	1470	1471	1476	rVB	1277	1693	0.15%	0.015%
29	12.206	1546	1552	1558	rBV3	2392	4850	0.44%	0.042%
30	12.917	1668	1673	1674	rBV3	2094	3101	0.28%	0.027%
31	13.546	1776	1780	1785	rVB2	2437	3349	0.30%	0.029%
32	13.698	1800	1806	1815	rBV	68938	115714	10.46%	0.999%
33	13.875	1830	1836	1843	rBV	364801	514242	46.48%	4.440%
34	14.162	1877	1885	1892	rVB	321655	509437	46.04%	4.399%
35	14.468	1930	1937	1945	rBV	293349	436480	39.45%	3.769%
36	14.668	1966	1971	1982	rBV2	35836	66953	6.05%	0.578%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091125.MA.M
 Title : SVOA CALIBRATION

37	15.009	2022	2029	2036	rBV	308881	487630	44.07%	4.210%
38	15.155	2050	2054	2059	rBV2	9377	14598	1.32%	0.126%
39	15.220	2063	2065	2068	rVB	1971	1700	0.15%	0.015%
40	15.273	2068	2074	2077	rBV4	2784	4757	0.43%	0.041%
41	15.461	2099	2106	2114	rVB	465402	681882	61.63%	5.887%
42	15.573	2119	2125	2133	rBV	215145	315635	28.53%	2.725%
43	15.825	2162	2168	2174	rVB	34791	50394	4.55%	0.435%
44	16.013	2196	2200	2201	rBV4	2926	2274	0.21%	0.020%
45	16.101	2213	2215	2219	rVB2	1337	1657	0.15%	0.014%
46	16.137	2219	2221	2225	rBV	1339	1556	0.14%	0.013%
47	16.331	2248	2254	2256	rBV	3470	4285	0.39%	0.037%
48	16.460	2272	2276	2278	rBV	1496	2232	0.20%	0.019%
49	16.489	2278	2281	2284	rVB5	3192	3474	0.31%	0.030%
50	16.618	2299	2303	2309	rBV4	14160	20116	1.82%	0.174%
51	16.777	2326	2330	2331	rBV2	1175	1496	0.14%	0.013%
52	16.971	2358	2363	2364	rBV	1257	1837	0.17%	0.016%
53	17.030	2369	2373	2374	rBV3	1476	1542	0.14%	0.013%
54	17.130	2386	2390	2394	rBV3	3325	4766	0.43%	0.041%
55	17.206	2395	2403	2409	rVV2	258437	375785	33.96%	3.245%
56	17.265	2411	2413	2414	rBV	2734	2643	0.24%	0.023%
57	17.306	2414	2420	2429	rVB2	575838	850865	76.90%	7.347%
58	17.382	2431	2433	2438	rVB3	5529	5742	0.52%	0.050%
59	17.447	2442	2444	2448	rVB2	1447	1776	0.16%	0.015%
60	17.488	2448	2451	2452	rBV2	1396	1488	0.13%	0.013%
61	17.600	2465	2470	2476	rBV5	4757	9718	0.88%	0.084%
62	17.717	2486	2490	2497	rVB4	5125	9452	0.85%	0.082%
63	17.794	2502	2503	2507	rBV	1396	1622	0.15%	0.014%
64	17.976	2528	2534	2540	rBV3	22593	35164	3.18%	0.304%
65	18.105	2551	2556	2564	rBV2	79489	114952	10.39%	0.993%
66	18.187	2564	2570	2576	rVB	175133	260850	23.58%	2.252%
67	18.405	2606	2607	2610	rVB3	2485	1777	0.16%	0.015%
68	18.469	2612	2618	2625	rBV	77419	111756	10.10%	0.965%
69	18.598	2638	2640	2642	rBV2	2127	1761	0.16%	0.015%
70	18.775	2667	2670	2674	rVB5	3176	3315	0.30%	0.029%
71	18.927	2690	2696	2709	rBV	259896	382377	34.56%	3.302%
72	19.086	2721	2723	2725	rBV2	2650	2820	0.25%	0.024%
73	19.157	2731	2735	2740	rBV3	9308	11701	1.06%	0.101%
74	19.233	2742	2748	2749	rBV2	10372	14028	1.27%	0.121%
75	19.257	2749	2752	2757	rVB6	9967	14727	1.33%	0.127%
76	19.380	2769	2773	2779	rBV4	23062	34673	3.13%	0.299%

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77	19.427	2779	2781	2783	rVV3	2597	2024	0.18%	0.017%
78	19.527	2794	2798	2803	rBV5	7049	12360	1.12%	0.107%
79	19.591	2803	2809	2817	rBV	734531	1047540	94.68%	9.045%
80	19.979	2873	2875	2878	rVB2	2716	1807	0.16%	0.016%
81	20.208	2912	2914	2917	rVB3	3766	3402	0.31%	0.029%
82	20.455	2954	2956	2960	rBV5	3860	3759	0.34%	0.032%
83	20.749	3004	3006	3010	rBV5	3174	3858	0.35%	0.033%
84	21.107	3060	3067	3072	rBV6	27746	68395	6.18%	0.591%
85	21.266	3091	3094	3097	rBV4	4944	7021	0.63%	0.061%
86	21.336	3104	3106	3112	rVB4	12449	15822	1.43%	0.137%
87	21.430	3116	3122	3129	rBV	302682	480125	43.39%	4.145%
88	21.636	3155	3157	3160	rBV4	4511	4308	0.39%	0.037%
89	21.848	3191	3193	3195	rBV3	4396	3285	0.30%	0.028%
90	22.306	3269	3271	3275	rBV4	4861	5389	0.49%	0.047%
91	23.023	3387	3393	3399	rVB	130084	225236	20.36%	1.945%
92	23.234	3422	3429	3439	rBV3	47450	125650	11.36%	1.085%
93	23.587	3487	3489	3496	rVB6	11358	16173	1.46%	0.140%
94	24.233	3589	3599	3609	rVB	441316	1106457	100.00%	9.553%
95	24.433	3624	3633	3645	rVB3	190700	525153	47.46%	4.534%
96	25.138	3751	3753	3754	rBV2	4343	2412	0.22%	0.021%
97	27.318	4122	4124	4125	rBV2	3914	2498	0.23%	0.022%
98	27.371	4132	4133	4135	rBV	3336	2666	0.24%	0.023%
99	30.749	4706	4708	4709	rBV	3942	2164	0.20%	0.019%
100	31.560	4844	4846	4848	rBV3	3665	3687	0.33%	0.032%

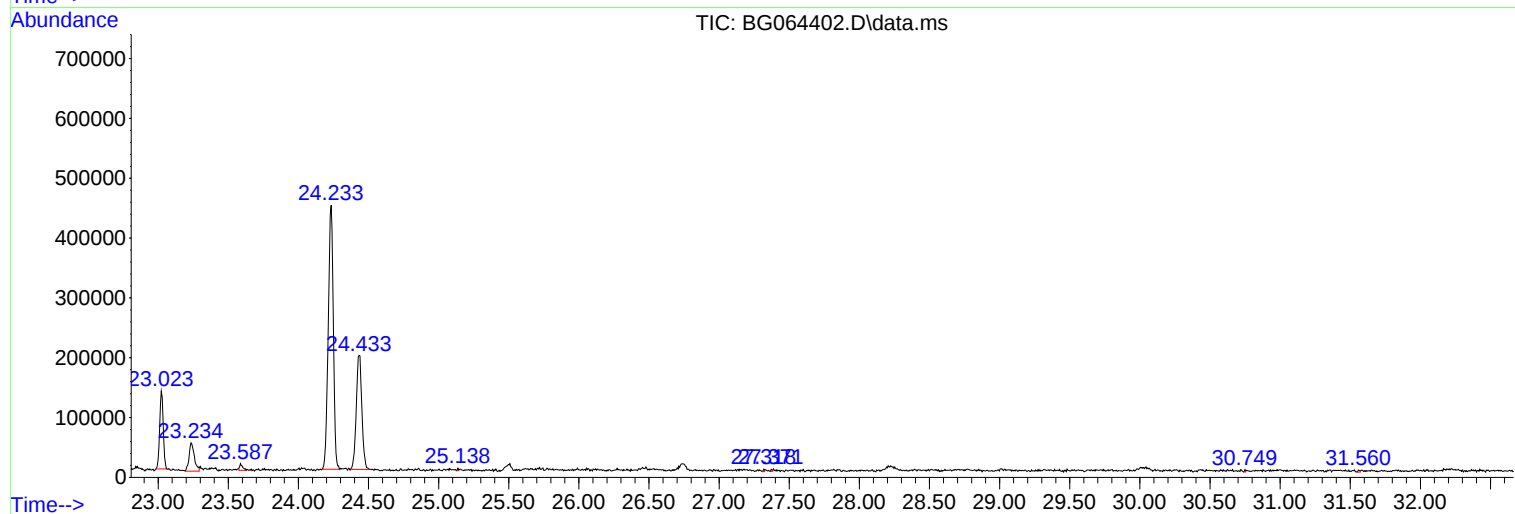
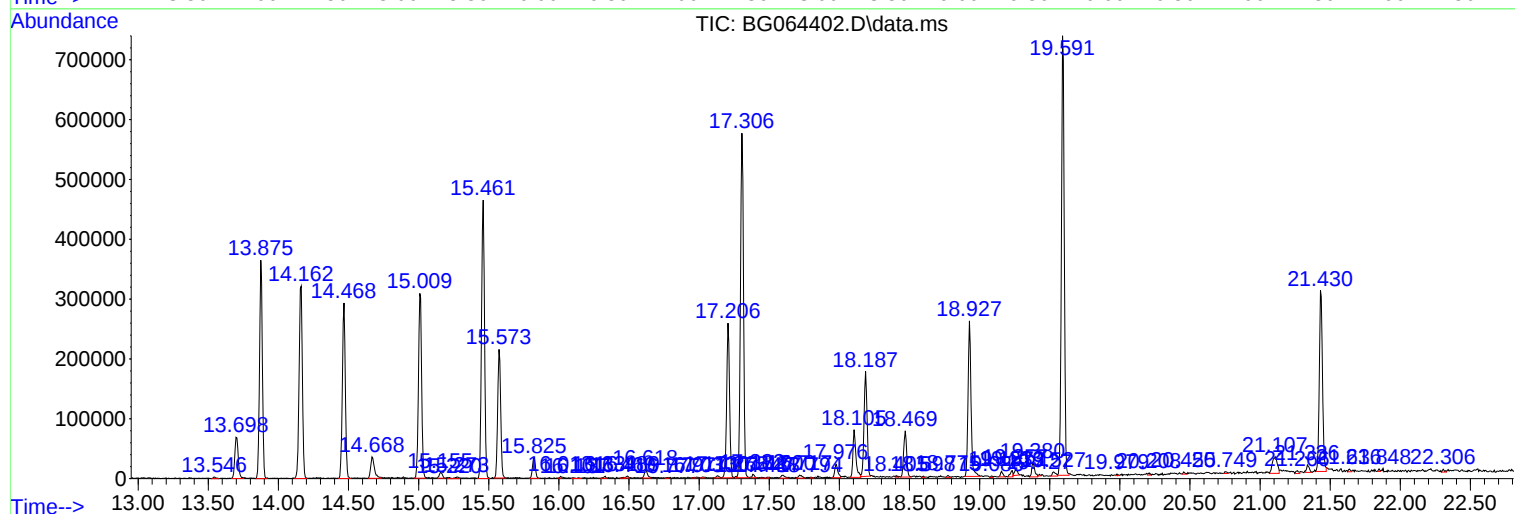
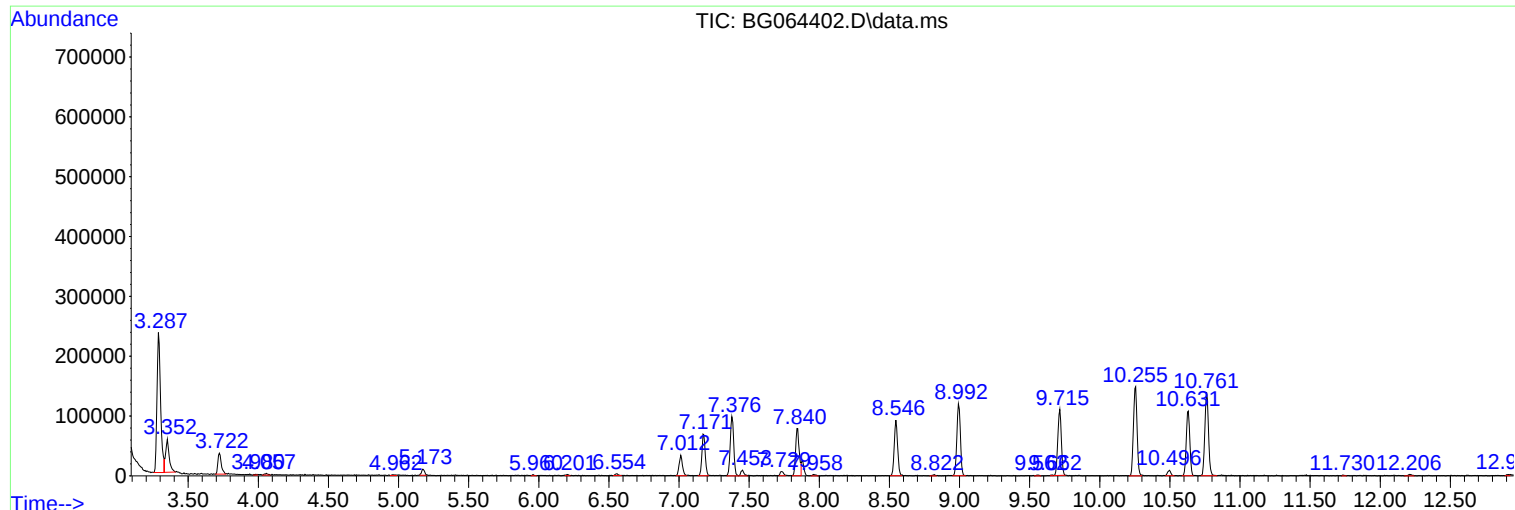
Sum of corrected areas: 11581907

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PMW-2(20250911)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091125.MA.M
Quant Title : SVOA CALIBRATION

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



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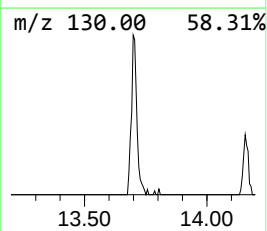
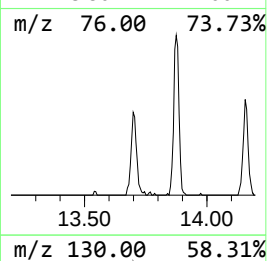
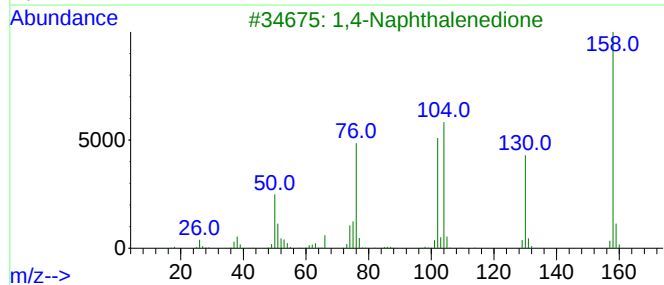
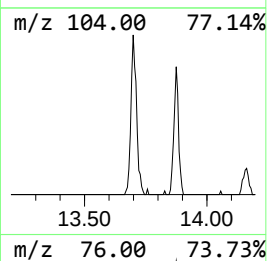
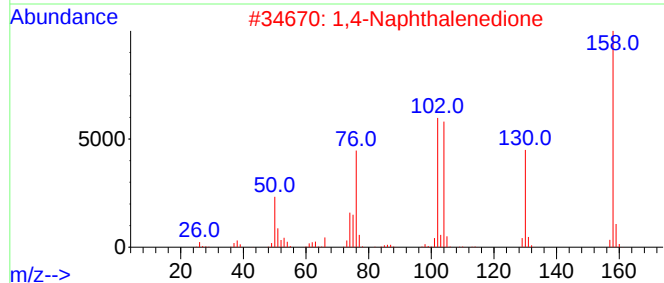
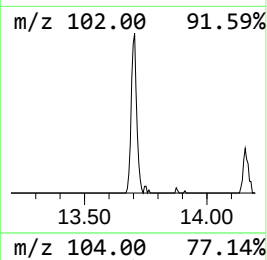
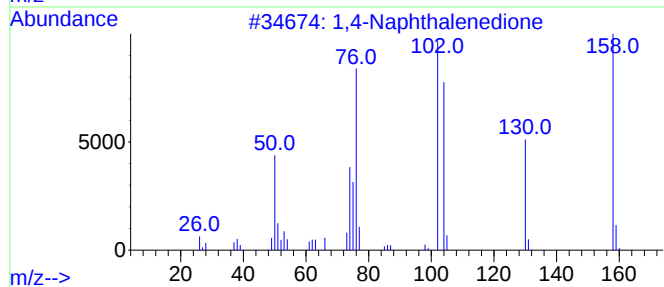
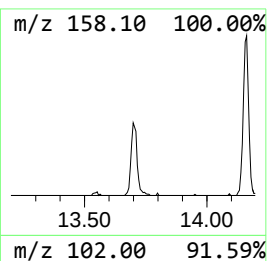
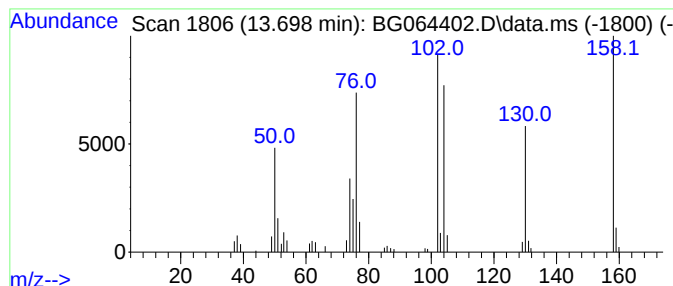
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091125.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,4-Naphthalenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.698	5.30 ng/ul	115714	Acenaphthene-d10	14.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Naphthalenedione			158	C10H6O2	000130-15-4	95
2	1,4-Naphthalenedione			158	C10H6O2	000130-15-4	94
3	1,4-Naphthalenedione			158	C10H6O2	000130-15-4	94
4	1,4-Naphthalenedione			158	C10H6O2	000130-15-4	76
5	Quinoxaline			130	C8H6N2	000091-19-0	58



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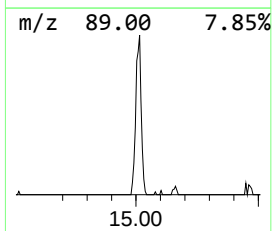
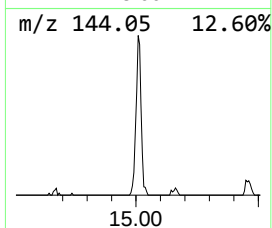
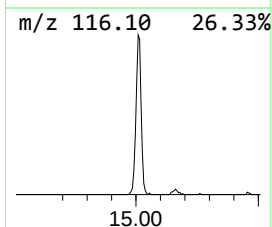
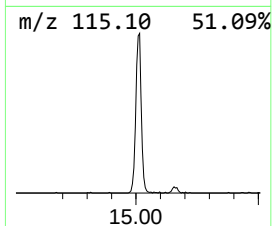
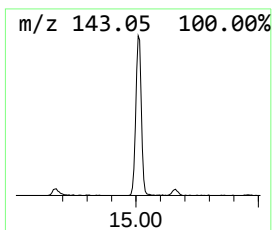
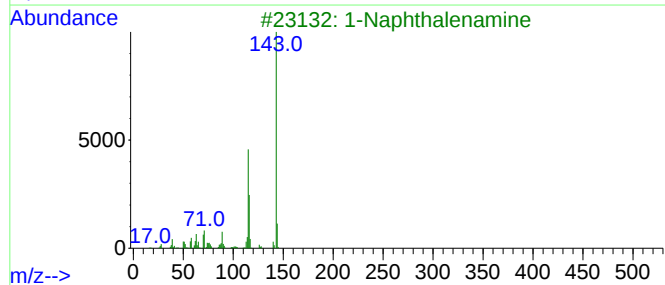
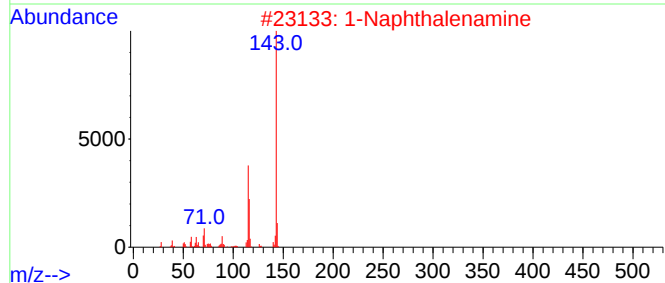
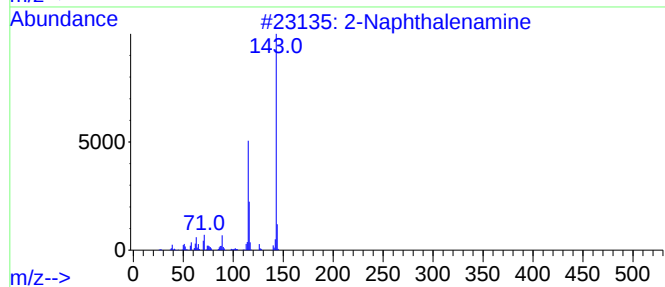
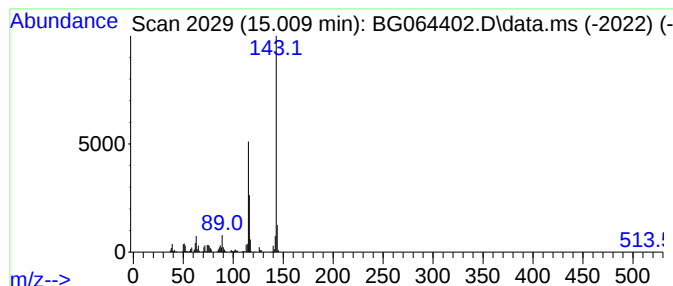
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091125.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Naphthalenamine Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.009	22.34 ng/ul	487630	Acenaphthene-d10	14.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenamine			143	C10H9N	000091-59-8	96
2	1-Naphthalenamine			143	C10H9N	000134-32-7	95
3	1-Naphthalenamine			143	C10H9N	000134-32-7	93
4	1H-Pyrrole, 2-phenyl-			143	C10H9N	003042-22-6	91
5	2-Naphthalenamine			143	C10H9N	000091-59-8	91



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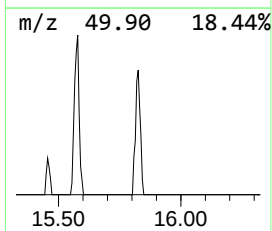
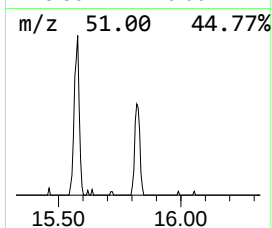
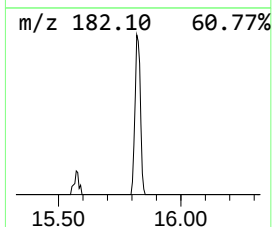
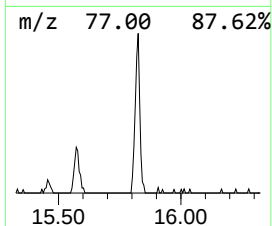
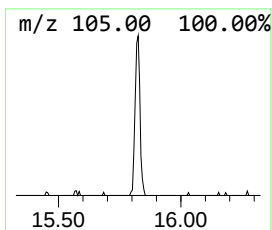
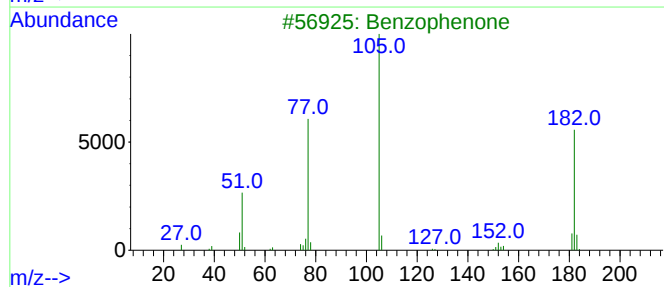
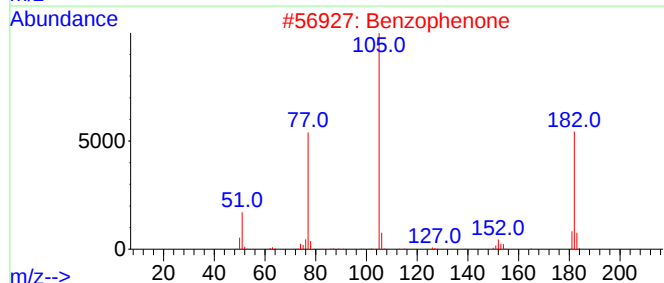
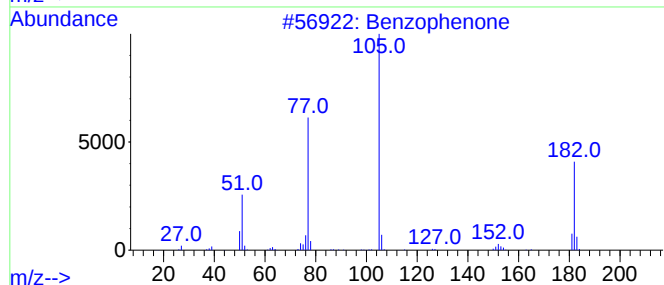
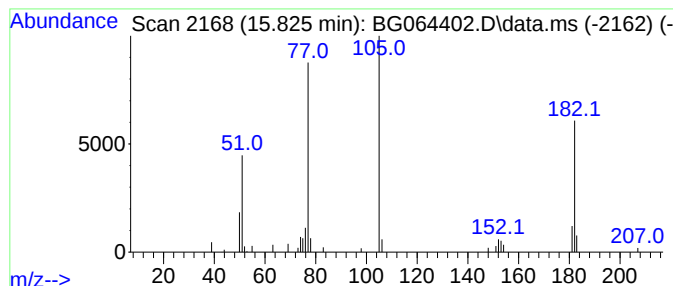
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzophenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.825	2.31 ng/ul	50394	Acenaphthene-d10	14.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	94
2			Benzophenone	182	C13H10O	000119-61-9	91
3			Benzophenone	182	C13H10O	000119-61-9	91
4			Benzophenone	182	C13H10O	000119-61-9	90
5			Benzophenone	182	C13H10O	000119-61-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091625\
Data File : BG064402.D
Acq On : 17 Sep 2025 2:58
Operator : RC/JU
Sample : Q3084-18
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PMW-2(20250911)

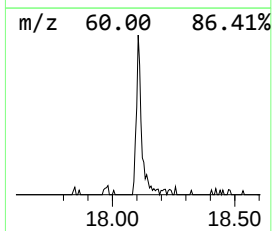
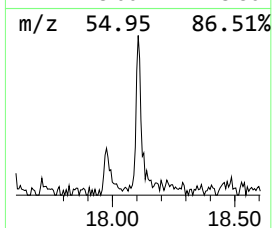
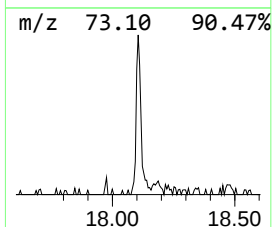
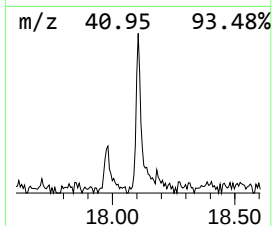
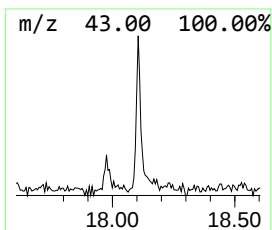
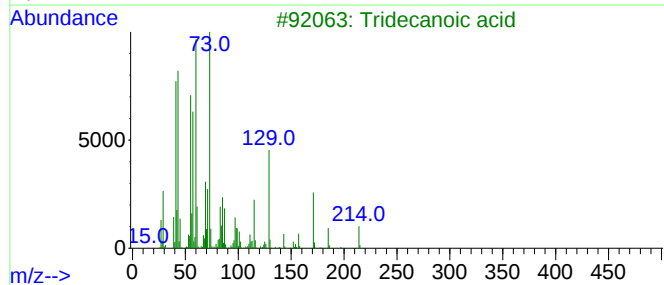
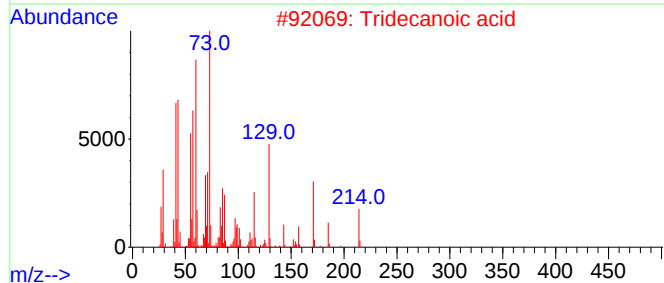
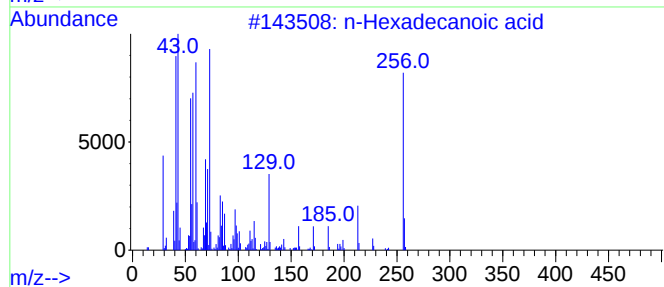
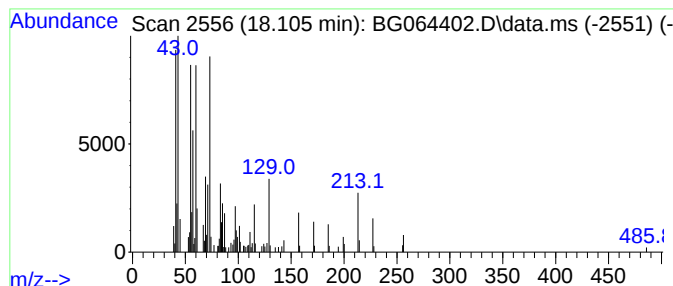
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.105	6.12 ng/ul	114952	Phenanthrene-d10	17.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2			Tridecanoic acid	214	C13H26O2	000638-53-9	96
3			Tridecanoic acid	214	C13H26O2	000638-53-9	95
4			Tridecanoic acid	214	C13H26O2	000638-53-9	89
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091625\
Data File : BG064402.D
Acq On : 17 Sep 2025 2:58
Operator : RC/JU
Sample : Q3084-18
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PMW-2(20250911)

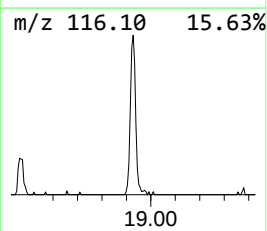
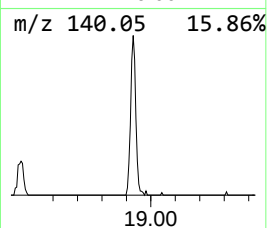
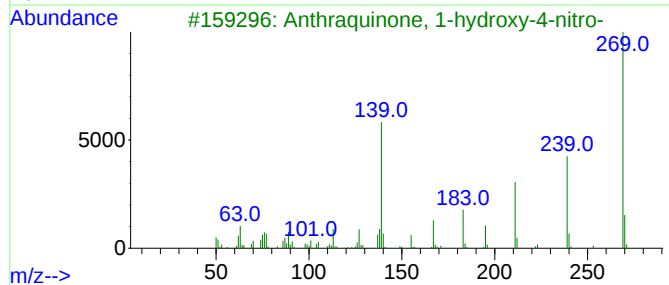
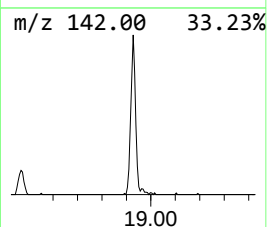
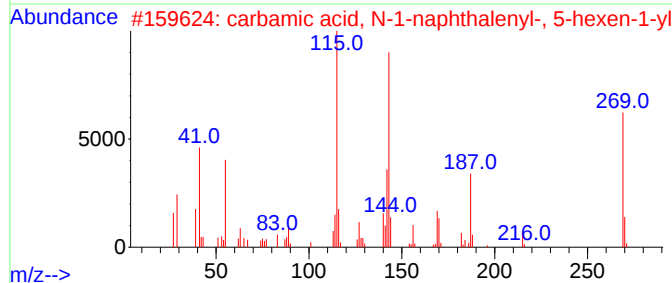
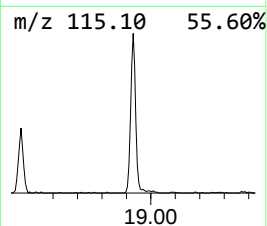
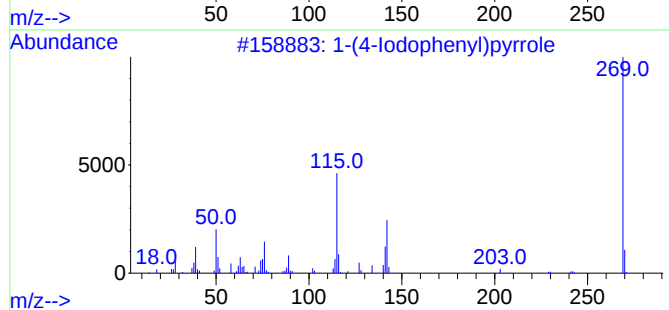
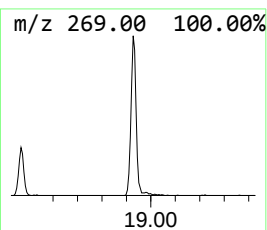
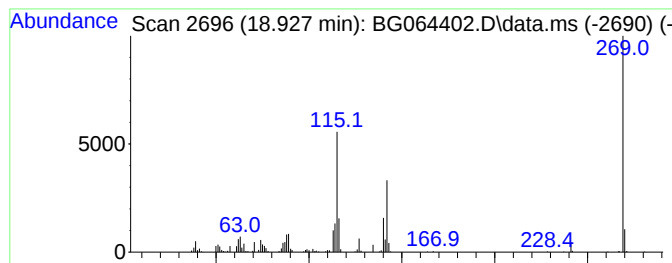
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1-(4-Iodophenyl)pyrrole Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.927	20.35 ng/ul	382377	Phenanthrene-d10	17.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-(4-Iodophenyl)pyrrole	269	C10H8IN	092636-36-7	90
2			carbamic acid, N-1-naphthalenyl-...	269	C17H19NO2	1010396-97-3	42
3			Anthraquinone, 1-hydroxy-4-nitro-	269	C14H7NO5	000081-65-2	38
4			Malononitrile, phenyl-	142	C9H6N2	003041-40-5	38
5			Propanamide, N-(4-methoxyphenyl)...	269	C10H8F5NO2	1000307-33-2	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091625\
Data File : BG064402.D
Acq On : 17 Sep 2025 2:58
Operator : RC/JU
Sample : Q3084-18
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PMW-2(20250911)

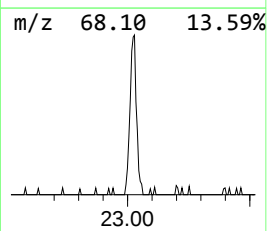
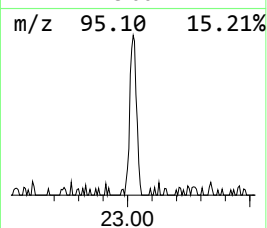
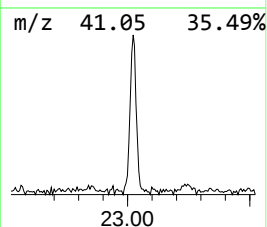
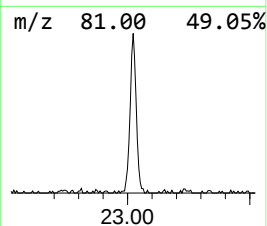
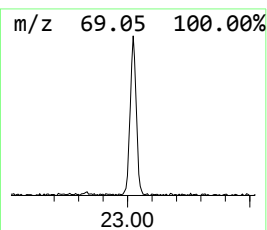
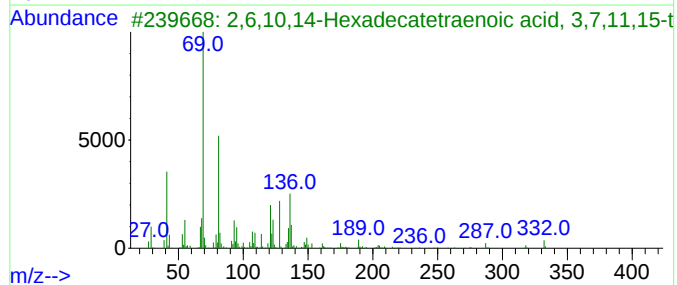
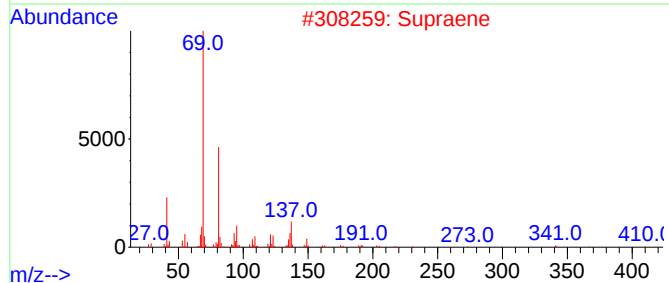
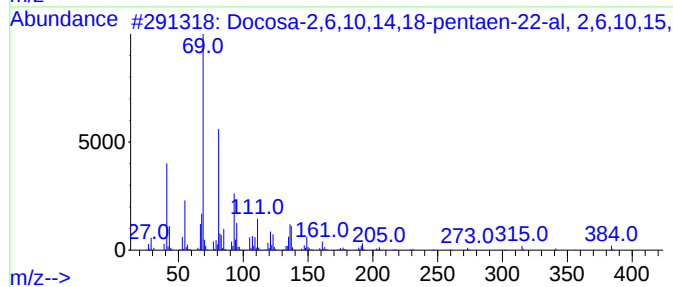
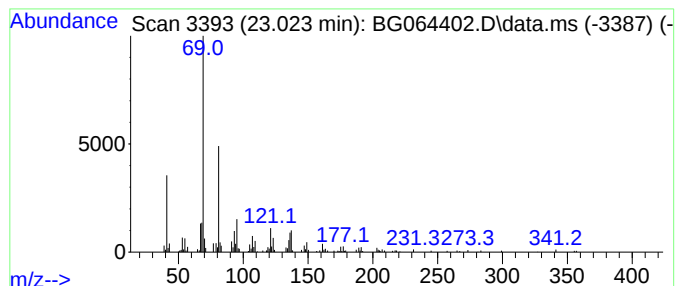
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Docosa-2,6,10,14,18-pentaen... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.023	8.58 ng/ul	225236	Perylene-d12	24.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	91
2			Supraene	410	C30H50	007683-64-9	80
3			2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	72
4			1,6-Octadiene, 3,5-dimethyl-, tr...	138	C10H18	074630-87-8	50
5			2-propenoic acid, 2-methyl-, 4-f...	190	C11H10O3	1000400-21-0	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091625\
Data File : BG064402.D
Acq On : 17 Sep 2025 2:58
Operator : RC/JU
Sample : Q3084-18
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PMW-2(20250911)

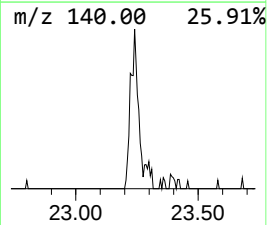
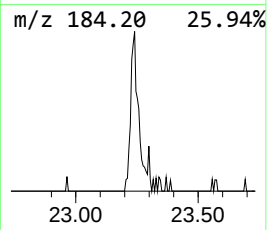
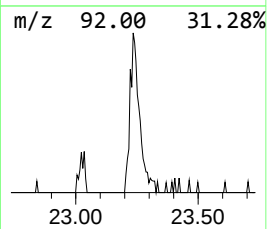
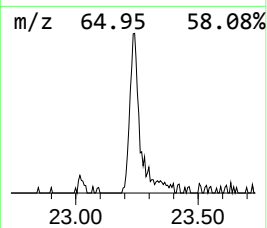
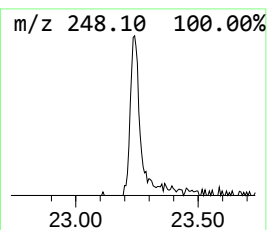
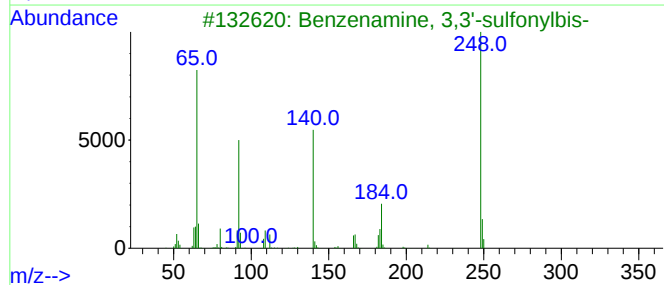
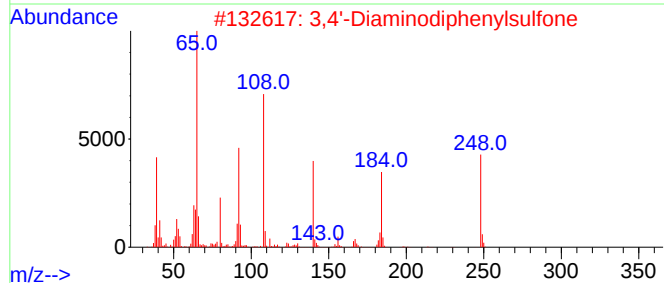
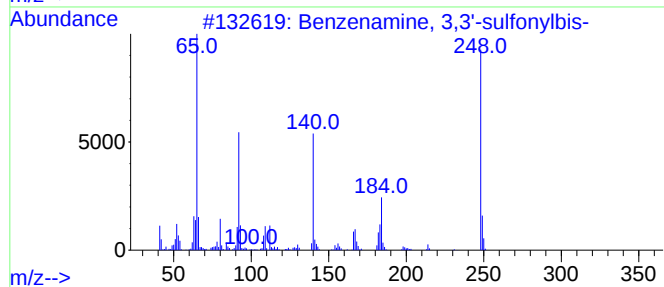
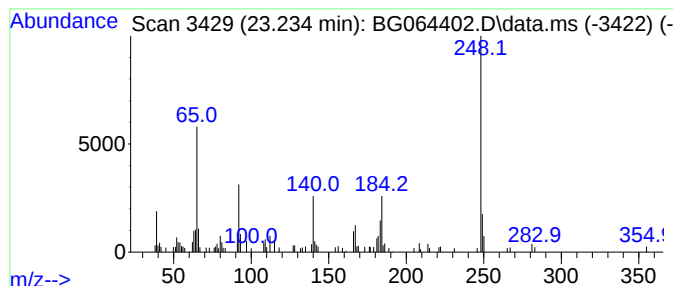
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzenamine, 3,3'-sulfonylbis- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.234	4.79 ng/ul	125650	Perylene-d12	24.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 3,3'-sulfonylbis-	248	C12H12N2O2S	000599-61-1	95
2			3,4'-Diaminodiphenylsulfone	248	C12H12N2O2S	034262-32-3	90
3			Benzenamine, 3,3'-sulfonylbis-	248	C12H12N2O2S	000599-61-1	72
4			Benzenamine, 3,3'-sulfonylbis-	248	C12H12N2O2S	000599-61-1	62
5			5,6,8,9-Tetrahydrobenz(a)anthrac...	248	C18H16O	001470-04-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091625\
Data File : BG064402.D
Acq On : 17 Sep 2025 2:58
Operator : RC/JU
Sample : Q3084-18
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PMW-2(20250911)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091125.MA.M
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
1,4-Naphthalene...	13.698	5.3	ng/ul	115714	3	14.468	436480	20.0
2-Naphthalenamine	15.009	22.3	ng/ul	487630	3	14.468	436480	20.0
Benzophenone	15.825	2.3	ng/ul	50394	3	14.468	436480	20.0
n-Hexadecanoic ...	18.105	6.1	ng/ul	114952	4	17.206	375785	20.0
1-(4-Iodophenyl...	18.927	20.4	ng/ul	382377	4	17.206	375785	20.0
Docosa-2,6,10,1...	23.023	8.6	ng/ul	225236	6	24.433	525153	20.0
Benzenamine, 3,...	23.234	4.8	ng/ul	125650	6	24.433	525153	20.0