

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
 Data File : BN015921.D
 Acq On : 18 Aug 2021 05:03
 Operator : CG/JU
 Sample : M3364-15
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBK99

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-BN081621MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN015921.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	38	42	50	rVB	75072	115335	1.39%	0.093%
2	3.758	112	114	127	rBV2	98134	179337	2.16%	0.145%
3	3.981	147	152	158	rVB2	29458	50098	0.60%	0.041%
4	7.058	668	675	688	rVB	308639	563612	6.78%	0.456%
5	7.228	697	704	713	rBV	931830	1585485	19.09%	1.283%
6	7.428	731	738	746	rBV	1051186	1766439	21.26%	1.430%
7	7.546	753	758	763	rVB3	28396	47815	0.58%	0.039%
8	7.899	811	818	827	rBV	981761	1677917	20.20%	1.358%
9	8.605	932	938	950	rVB	840223	1457029	17.54%	1.179%
10	9.063	1009	1016	1030	rVB	1304352	2270476	27.33%	1.838%
11	9.387	1067	1071	1077	rVV3	25841	51657	0.62%	0.042%
12	9.599	1102	1107	1113	rVB	28192	47700	0.57%	0.039%
13	9.787	1131	1139	1147	rBV	1061961	1806510	21.75%	1.462%
14	10.110	1188	1194	1201	rBV3	58576	112758	1.36%	0.091%
15	10.328	1221	1231	1242	rBV	1541796	2619747	31.54%	2.120%
16	10.710	1286	1296	1300	rBV	1382422	2478907	29.84%	2.006%
17	10.763	1300	1305	1312	rVB	3462861	5917132	71.23%	4.789%
18	10.846	1312	1319	1324	rBV	1240284	2431290	29.27%	1.968%
19	11.499	1423	1430	1434	rBV7	23005	53494	0.64%	0.043%
20	11.634	1446	1453	1461	rVB2	27171	49533	0.60%	0.040%
21	11.822	1480	1485	1492	rVB	23986	43103	0.52%	0.035%
22	12.098	1527	1532	1538	rBV6	30499	57402	0.69%	0.046%
23	12.263	1555	1560	1571	rBV2	118252	251459	3.03%	0.204%
24	12.375	1574	1579	1584	rVB2	46650	80857	0.97%	0.065%
25	12.493	1591	1599	1603	rBV6	31529	63073	0.76%	0.051%
26	12.593	1605	1616	1627	rVB	2056659	3474963	41.83%	2.812%
27	13.381	1744	1750	1757	rBV	970559	1479260	17.81%	1.197%
28	13.581	1778	1784	1787	rBV2	149296	275434	3.32%	0.223%
29	13.728	1796	1809	1816	rBV2	273806	697713	8.40%	0.565%
30	13.869	1824	1833	1838	rVV2	519535	965680	11.63%	0.782%
31	13.957	1838	1848	1857	rVV	3101798	4900471	58.99%	3.966%
32	14.104	1867	1873	1877	rVV	232487	383968	4.62%	0.311%
33	14.140	1877	1879	1884	rVB	94853	116775	1.41%	0.095%
34	14.240	1889	1896	1908	rVB	3108906	5265357	63.39%	4.261%
35	14.551	1939	1949	1954	rBV	2050219	3398672	40.91%	2.751%
36	14.610	1954	1959	1966	rVV	1881755	2959985	35.63%	2.396%

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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 >
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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
 Title : SVOA CALIBRATION

37	14.675	1966	1970	1976	rVV	381290	572792	6.90%	0.464%
38	14.740	1976	1981	1992	rVB2	303724	558014	6.72%	0.452%
39	14.945	2010	2016	2024	rVV	2993099	4642049	55.88%	3.757%
40	15.022	2024	2029	2036	rVV3	58012	110524	1.33%	0.089%
41	15.163	2045	2053	2059	rBV3	52384	108773	1.31%	0.088%
42	15.216	2059	2062	2070	rBV5	42896	71730	0.86%	0.058%
43	15.339	2075	2083	2086	rBV6	38873	96721	1.16%	0.078%
44	15.469	2100	2105	2110	rBV5	58307	106099	1.28%	0.086%
45	15.545	2110	2118	2122	rVV	4128323	6297016	75.80%	5.096%
46	15.598	2122	2127	2132	rVV	1723273	2555134	30.76%	2.068%
47	15.657	2132	2137	2144	rVV	1350278	2039276	24.55%	1.650%
48	15.722	2144	2148	2152	rVV3	122122	226711	2.73%	0.183%
49	15.775	2155	2157	2161	rVV4	79976	138656	1.67%	0.112%
50	15.816	2161	2164	2167	rVV2	94428	140542	1.69%	0.114%
51	15.845	2167	2169	2173	rVV	73027	104722	1.26%	0.085%
52	15.892	2173	2177	2183	rVV	222426	322509	3.88%	0.261%
53	15.998	2187	2195	2197	rVV2	89862	161344	1.94%	0.131%
54	16.034	2197	2201	2211	rVV2	290524	602103	7.25%	0.487%
55	16.128	2213	2217	2223	rVB	41438	64761	0.78%	0.052%
56	16.263	2228	2240	2248	rBV	3046816	4719117	56.81%	3.819%
57	16.322	2248	2250	2256	rVV7	32089	68746	0.83%	0.056%
58	16.575	2286	2293	2296	rBV3	89862	176707	2.13%	0.143%
59	16.604	2296	2298	2302	rVV3	73735	96047	1.16%	0.078%
60	16.651	2303	2306	2312	rVV5	41295	75773	0.91%	0.061%
61	16.834	2328	2337	2341	rBV7	36969	60893	0.73%	0.049%
62	16.916	2344	2351	2353	rVV	946238	1374710	16.55%	1.113%
63	16.945	2353	2356	2368	rVV	2246642	3376469	40.65%	2.733%
64	17.075	2373	2378	2381	rVV5	64236	129857	1.56%	0.105%
65	17.116	2381	2385	2390	rVV	276208	439990	5.30%	0.356%
66	17.163	2390	2393	2399	rVB5	67296	101632	1.22%	0.082%
67	17.298	2409	2416	2420	rVV	2486755	3855457	46.41%	3.120%
68	17.339	2420	2423	2428	rVV	2605691	3671757	44.20%	2.972%
69	17.398	2428	2433	2444	rVV	5107298	7285876	87.71%	5.897%
70	17.492	2444	2449	2450	rVV5	35886	63309	0.76%	0.051%
71	17.522	2451	2454	2459	rVB4	46967	64713	0.78%	0.052%
72	17.698	2473	2484	2490	rBV	974686	1506679	18.14%	1.219%
73	17.798	2496	2501	2504	rBV3	41482	63045	0.76%	0.051%
74	17.916	2517	2521	2525	rVV6	25430	46750	0.56%	0.038%
75	17.975	2525	2531	2532	rVV5	30888	48454	0.58%	0.039%
76	18.004	2532	2536	2546	rVB4	82313	159962	1.93%	0.129%

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77	18.092	2546	2551	2559	rBV5	39142	75428	0.91%	0.061%
78	18.181	2560	2566	2570	rBV3	96518	192687	2.32%	0.156%
79	18.222	2570	2573	2577	rVV2	62823	90774	1.09%	0.073%
80	18.263	2577	2580	2587	rVB3	58142	93097	1.12%	0.075%
81	18.375	2594	2599	2603	rBV3	54992	103867	1.25%	0.084%
82	18.551	2627	2629	2639	rVB10	24194	55299	0.67%	0.045%
83	18.675	2646	2650	2653	rVV3	32352	50846	0.61%	0.041%
84	18.716	2653	2657	2662	rVB	195263	276416	3.33%	0.224%
85	19.092	2717	2721	2725	rBV4	38656	55924	0.67%	0.045%
86	19.227	2740	2744	2747	rBV2	50472	86899	1.05%	0.070%
87	19.322	2756	2760	2762	rBV	72431	99934	1.20%	0.081%
88	19.357	2762	2766	2771	rVV	175913	256396	3.09%	0.208%
89	19.416	2772	2776	2784	rVB2	97533	148059	1.78%	0.120%
90	19.692	2816	2823	2833	rBV2	6026890	8306911	100.00%	6.723%
91	19.780	2833	2838	2844	rVB	481418	657426	7.91%	0.532%
92	20.151	2896	2901	2906	rBV	126632	193185	2.33%	0.156%
93	20.692	2989	2993	2998	rBV	199451	235891	2.84%	0.191%
94	20.810	3009	3013	3018	rVB2	72583	102164	1.23%	0.083%
95	20.886	3021	3026	3032	rBV	153680	211275	2.54%	0.171%
96	21.192	3074	3078	3083	rBV2	221862	350996	4.23%	0.284%
97	21.398	3110	3113	3119	rBV	106972	135815	1.63%	0.110%
98	21.480	3122	3127	3131	rBV	2958431	4331068	52.14%	3.505%
99	23.751	3505	3513	3520	rBV	3760952	7667129	92.30%	6.205%
100	23.904	3532	3539	3549	rVB2	1808912	3778104	45.48%	3.058%

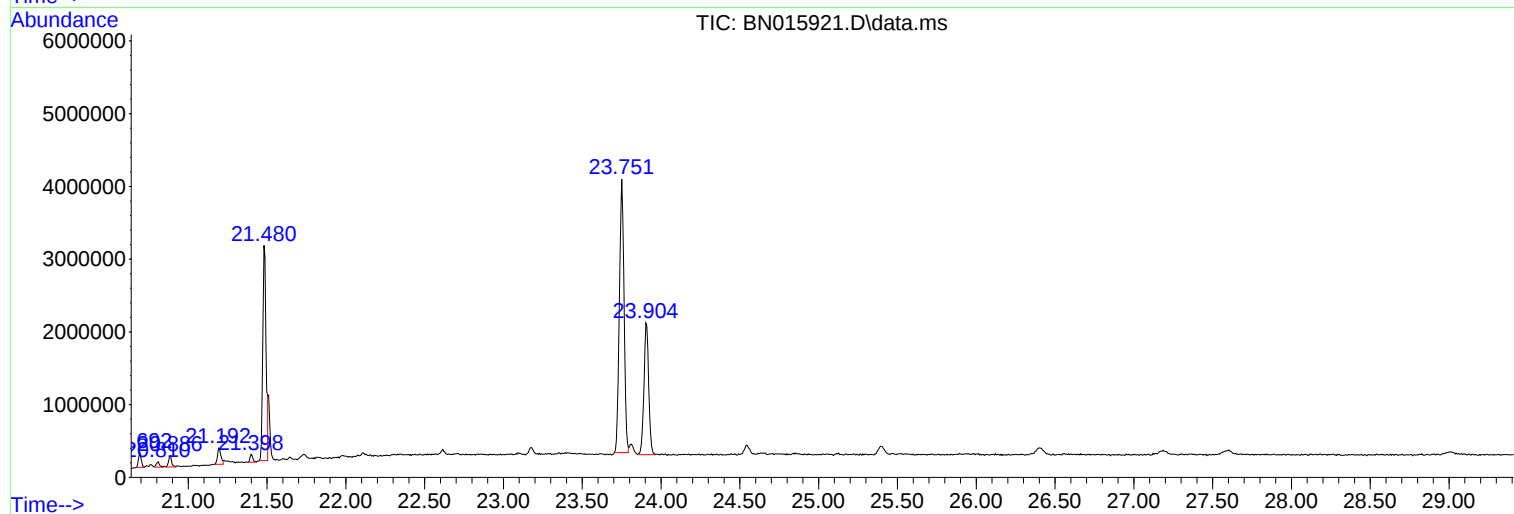
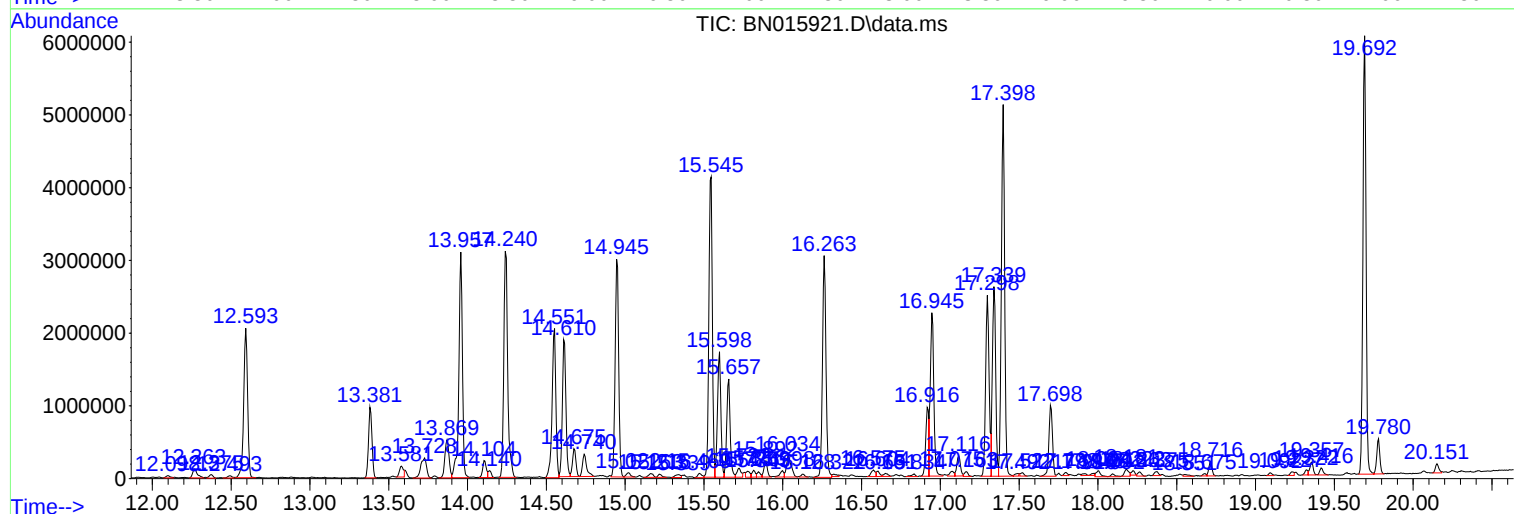
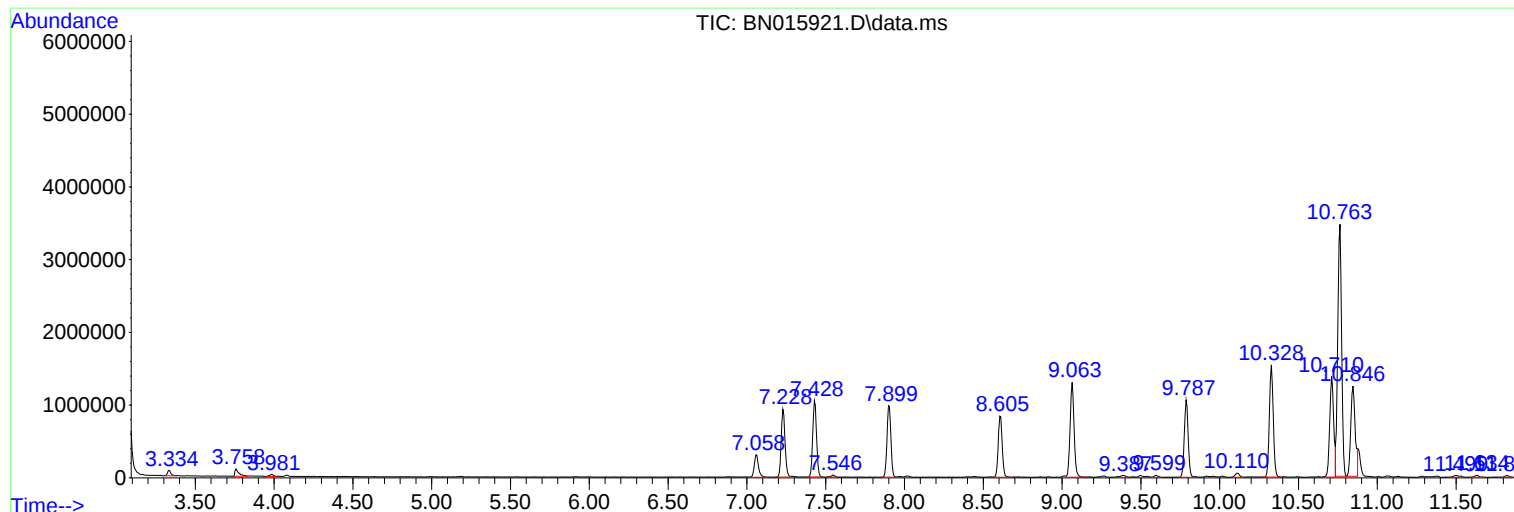
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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



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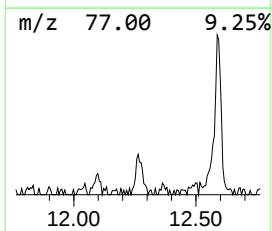
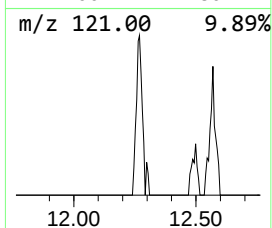
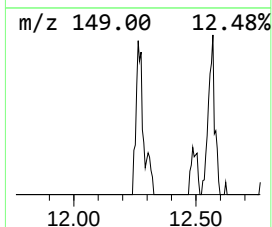
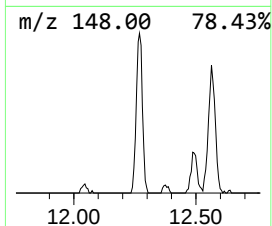
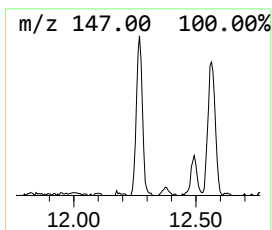
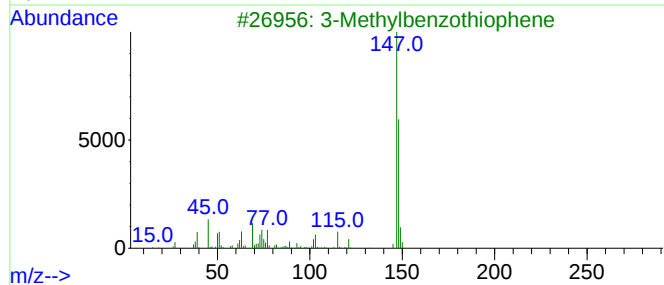
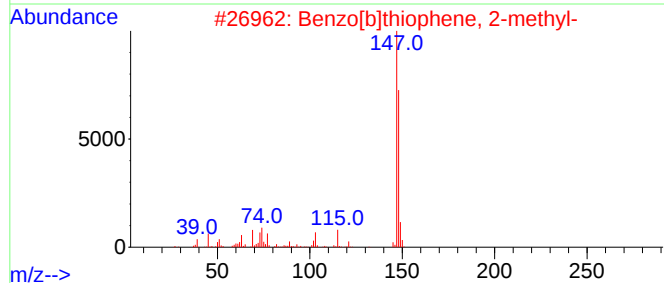
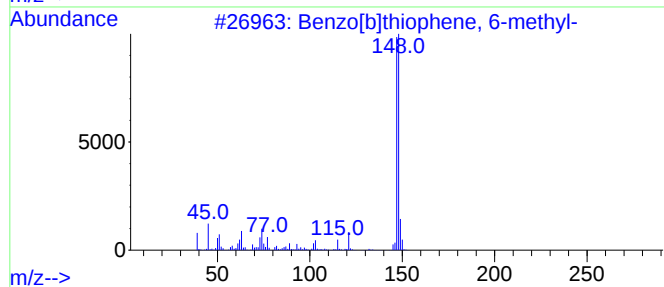
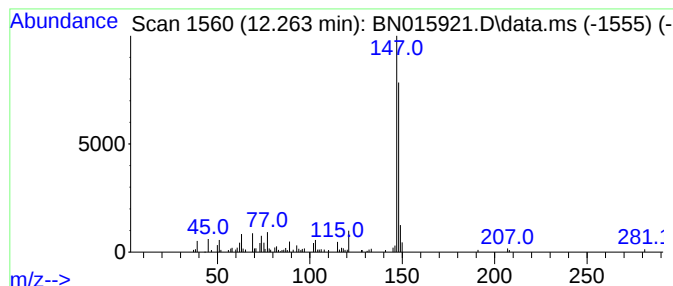
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzo[b]thiophene, 6-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.263	2.03 ng/ul	251459	Naphthalene-d8	10.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene, 6-methyl-	148	C9H8S	016587-47-6	93
2			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	90
3			3-Methylbenzothiophene	148	C9H8S	001455-18-1	87
4			Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	87
5			Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	83



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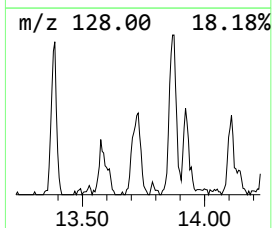
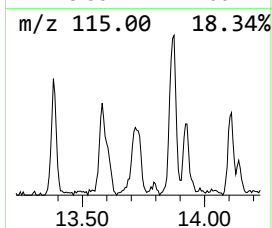
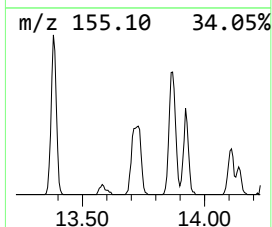
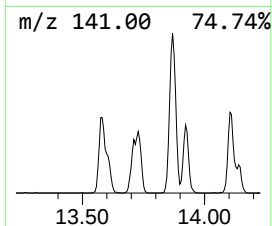
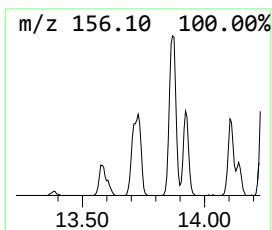
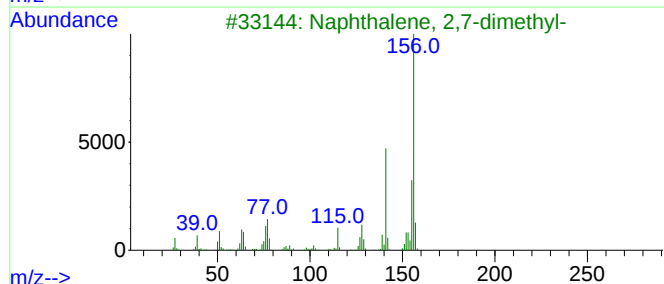
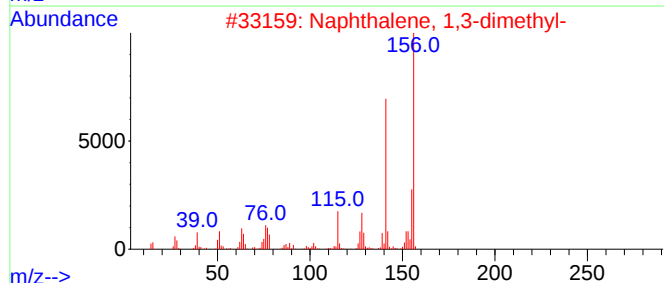
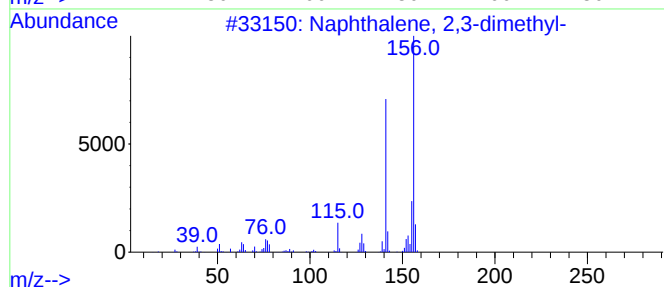
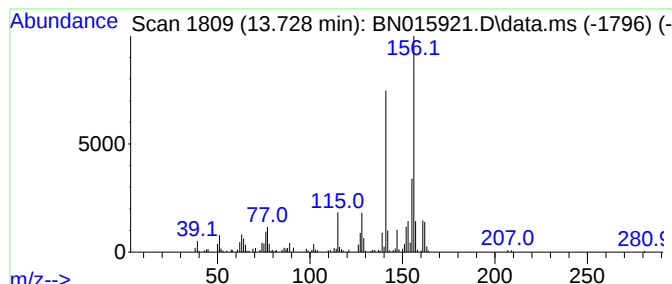
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Naphthalene, 2,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.728	4.11 ng/ul	697713	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96



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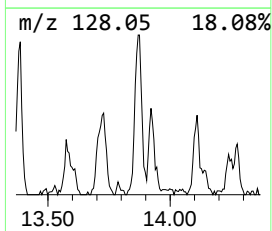
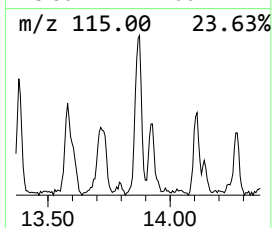
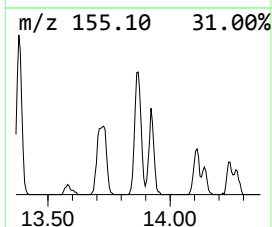
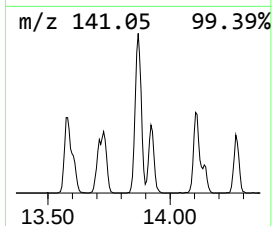
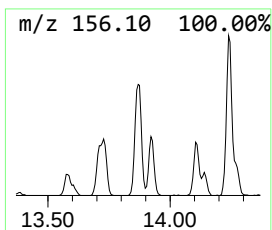
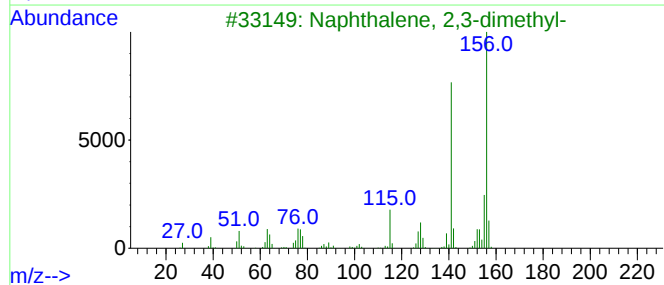
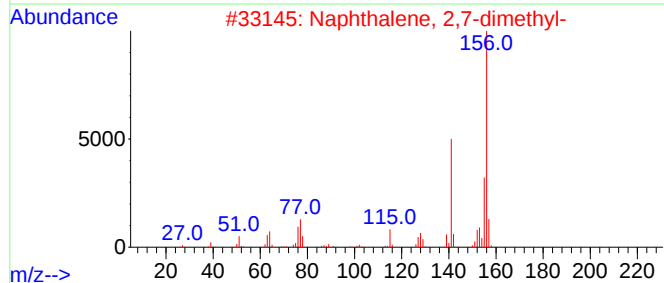
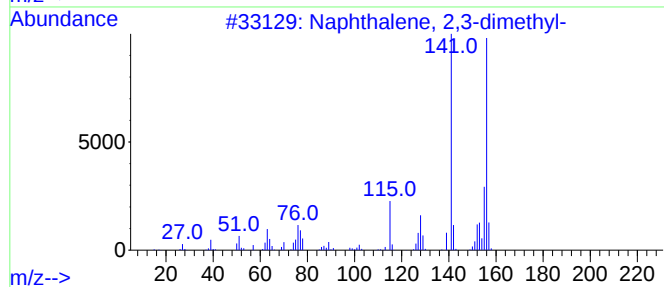
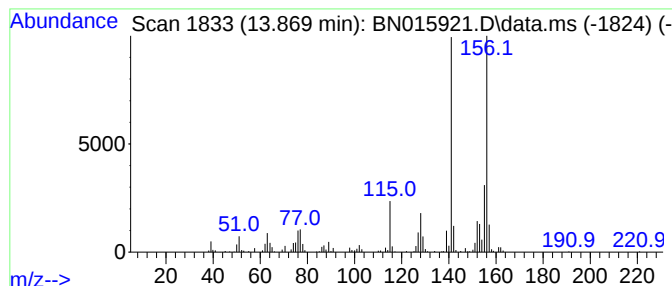
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Naphthalene, 2,7-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.869	5.68 ng/ul	965680	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015921.D
Acq On : 18 Aug 2021 05:03
Operator : CG/JU
Sample : M3364-15
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBK99

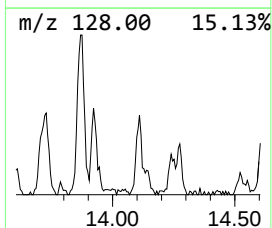
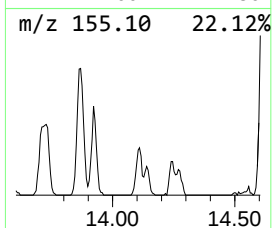
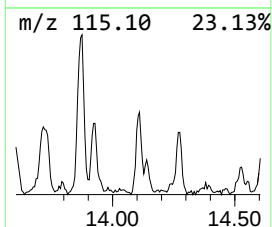
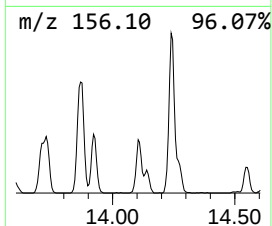
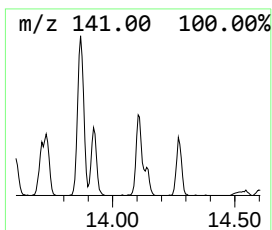
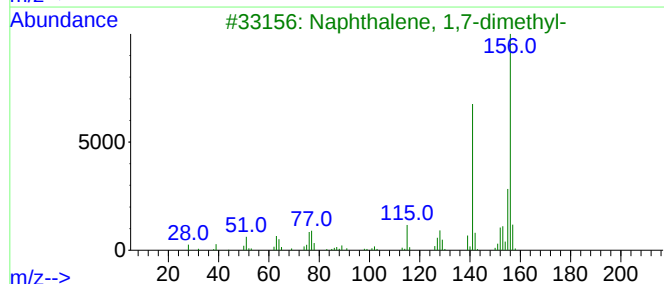
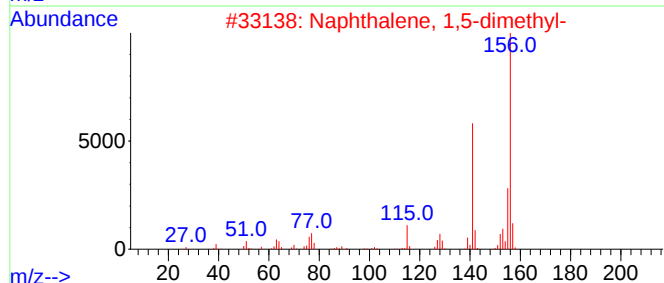
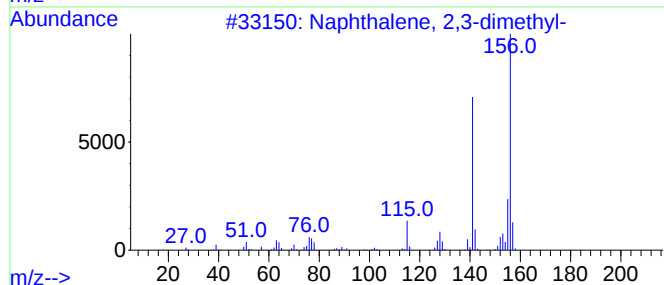
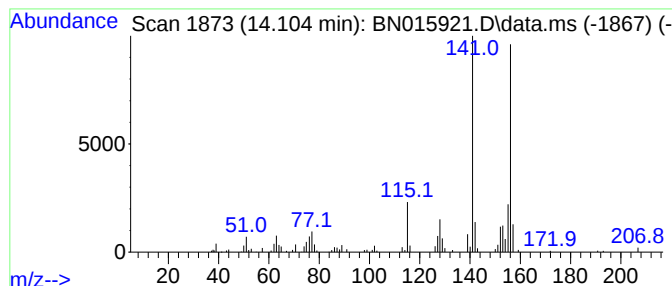
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Naphthalene, 1,5-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.104	2.26 ng/ul	383968	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
5			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015921.D
Acq On : 18 Aug 2021 05:03
Operator : CG/JU
Sample : M3364-15
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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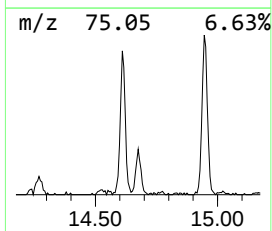
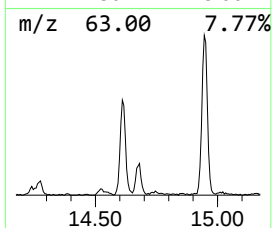
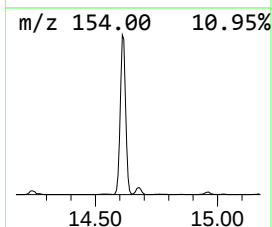
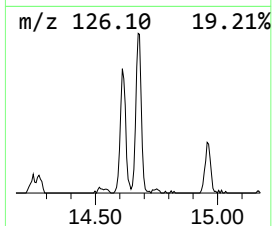
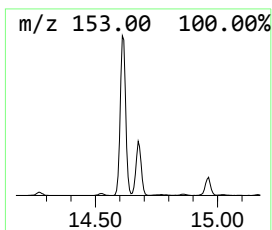
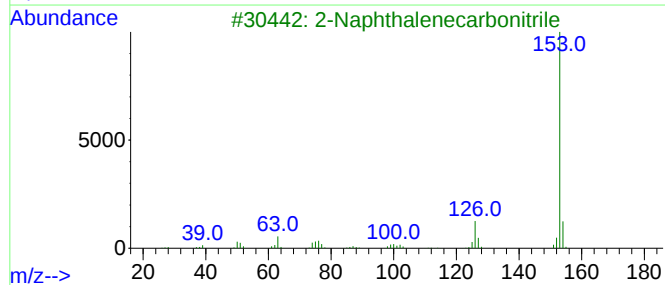
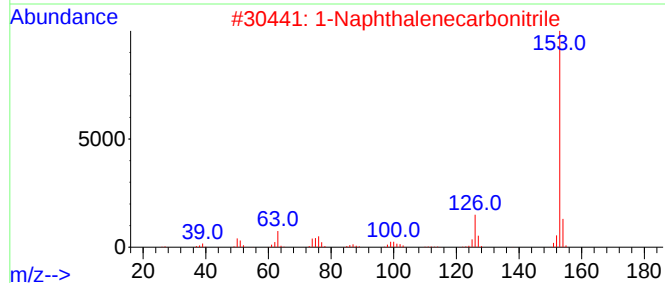
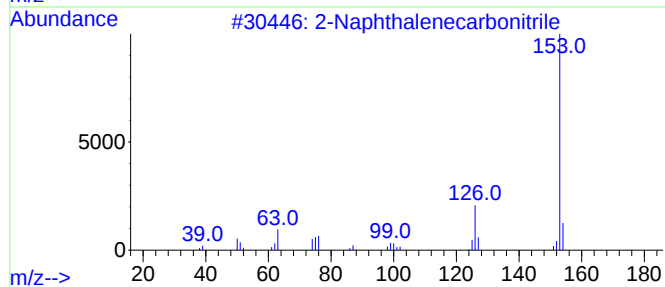
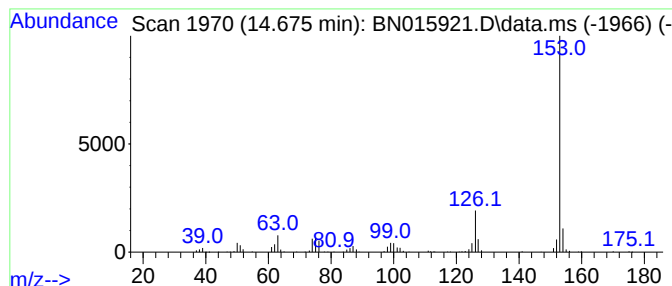
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2-Naphthalenecarbonitrile Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.675	3.37 ng/ul	572792	Acenaphthene-d10	14.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	97		
2	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	94		
3	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	91		
4	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	90		
5	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	90		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015921.D
Acq On : 18 Aug 2021 05:03
Operator : CG/JU
Sample : M3364-15
Misc :
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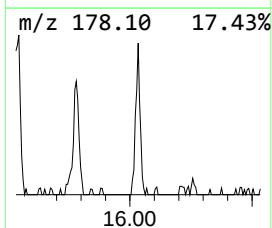
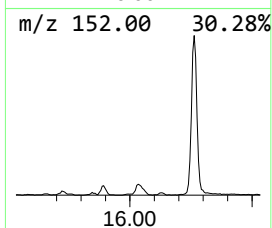
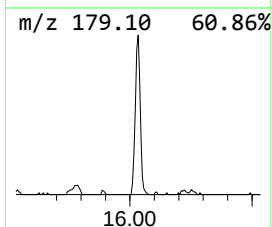
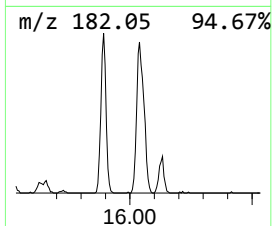
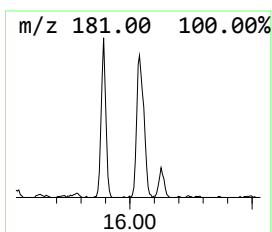
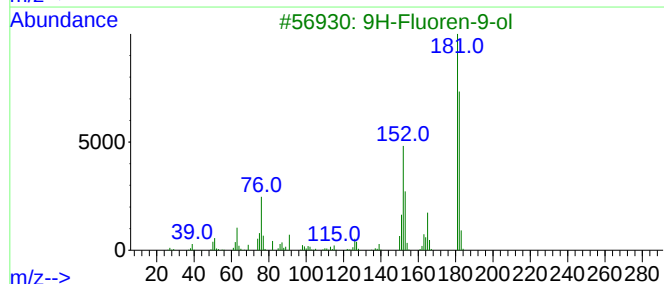
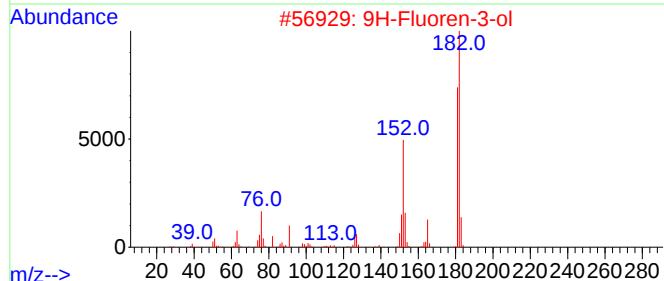
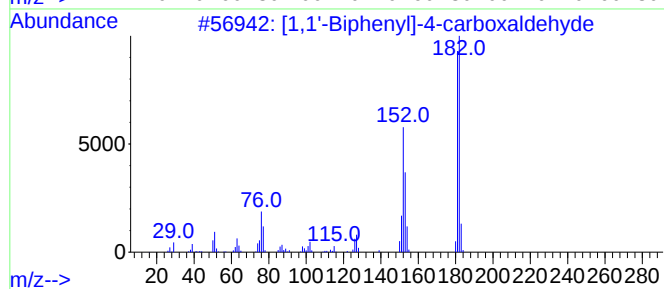
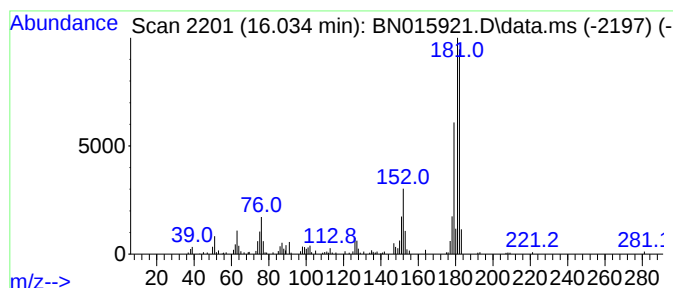
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 [1,1'-Biphenyl]-4-carboxald... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.034	3.12 ng/ul	602103	Phenanthrene-d10	17.298	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	86
2	9H-Fluoren-3-ol	182	C13H10O	006344-67-8	68
3	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	68
4	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	64
5	3-(1-Naphthyl)acrylaldehyde	182	C13H10O	001504-73-0	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015921.D
Acq On : 18 Aug 2021 05:03
Operator : CG/JU
Sample : M3364-15
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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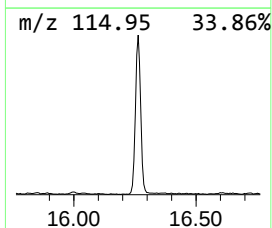
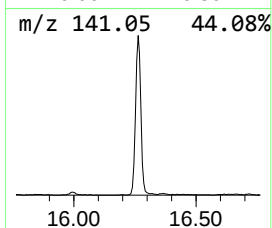
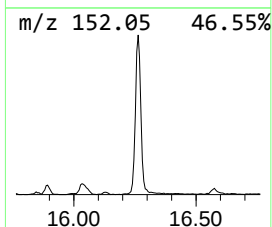
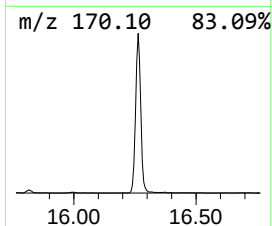
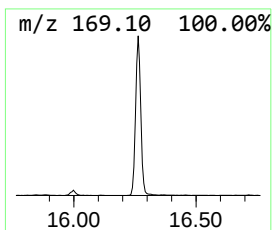
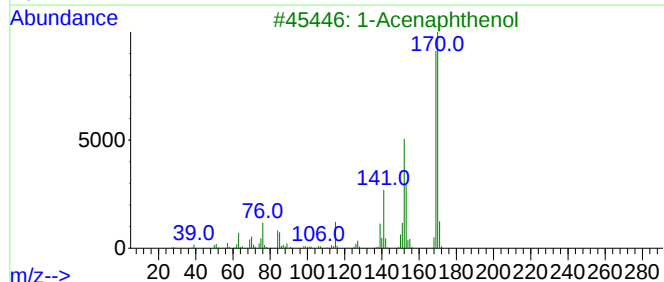
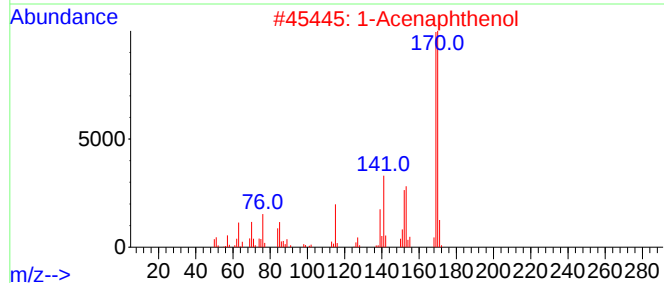
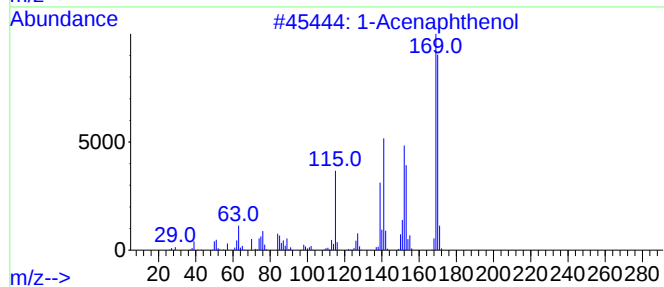
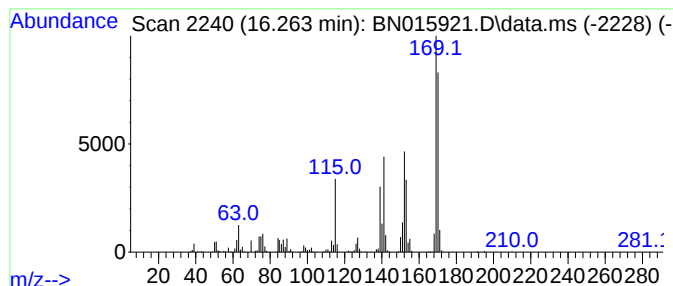
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Acenaphthenol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.263	24.48 ng/ul	4719120	Phenanthrene-d10	17.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Acenaphthenol	170	C12H10O	006306-07-6	98		
2	1-Acenaphthenol	170	C12H10O	006306-07-6	95		
3	1-Acenaphthenol	170	C12H10O	006306-07-6	70		
4	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	62		
5	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	53		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015921.D
Acq On : 18 Aug 2021 05:03
Operator : CG/JU
Sample : M3364-15
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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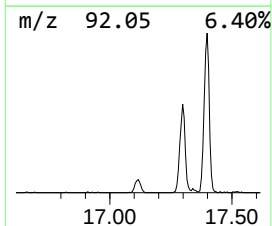
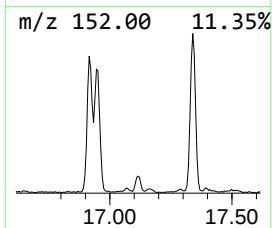
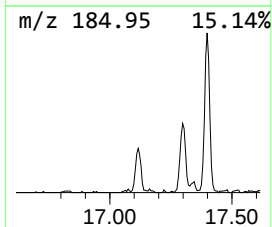
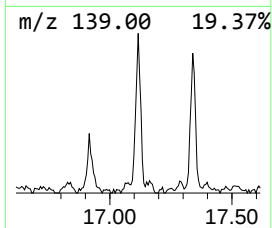
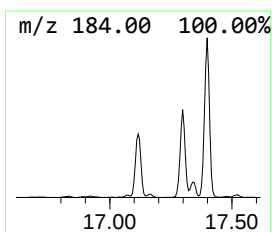
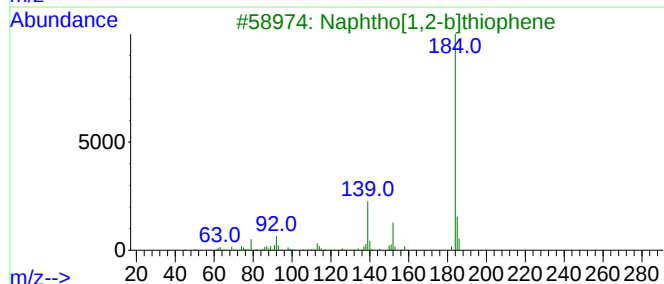
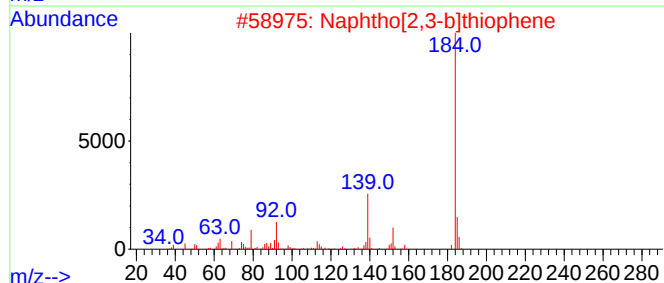
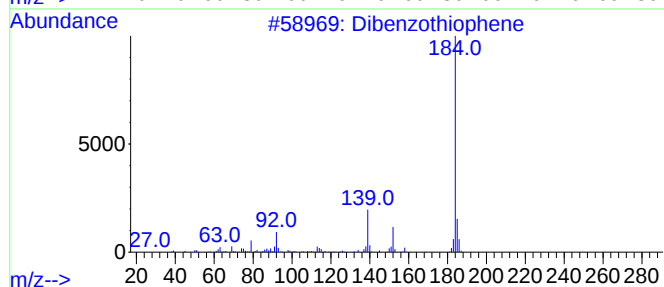
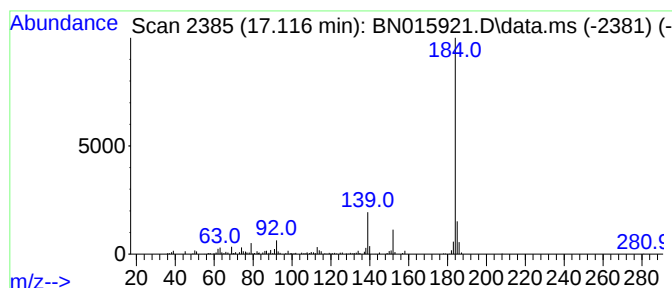
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Dibenzothiophene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.116	2.28 ng/ul	439990	Phenanthrene-d10	17.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	97
2			Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	97
3			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	96
4			Dibenzothiophene	184	C12H8S	000132-65-0	95
5			Dibenzothiophene	184	C12H8S	000132-65-0	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015921.D
Acq On : 18 Aug 2021 05:03
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Instrument :
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ClientSampleId :
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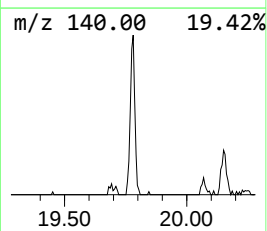
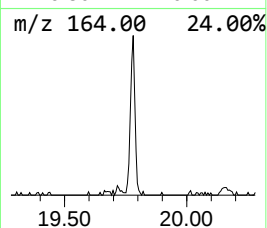
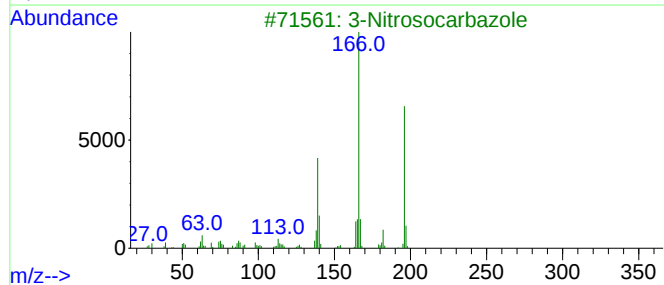
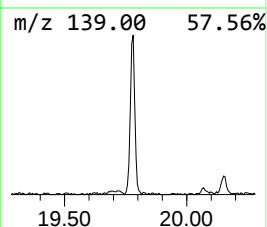
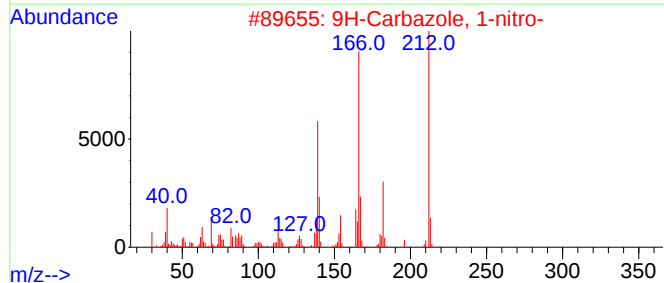
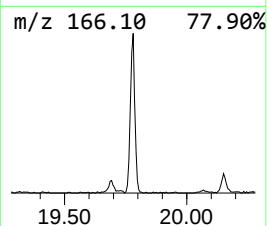
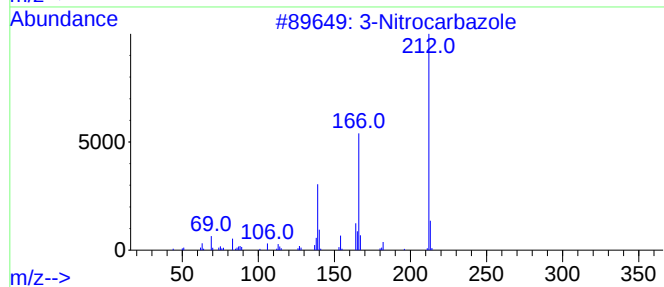
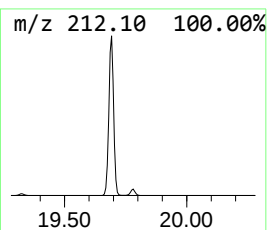
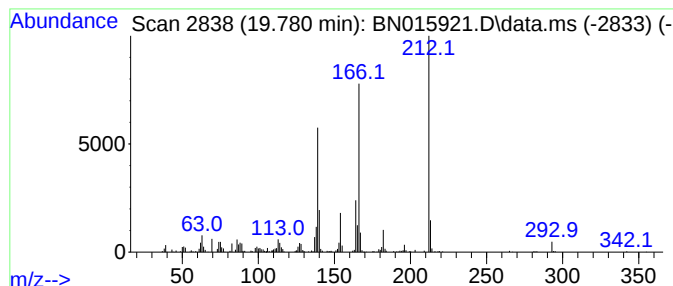
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 3-Nitrocarbazole Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.780	3.04 ng/ul	657426	Chrysene-d12	21.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Nitrocarbazole	212	C12H8N2O2	003077-85-8	96
2			9H-Carbazole, 1-nitro-	212	C12H8N2O2	031438-22-9	90
3			3-Nitrosocarbazole	196	C12H8N2O	040310-53-0	55
4			1H-Pyrrolo[3,2-h]quinoline-2-car...	212	C12H8N2O2	1000337-74-6	53
5			Benzo[b]benzofuran-2-carboxylic ...	212	C13H8O3	1010337-38-1	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015921.D
Acq On : 18 Aug 2021 05:03
Operator : CG/JU
Sample : M3364-15
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBK99

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
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
Benzo[b]thiophe...	12.263	2.0	ng/u1	251459	2	10.710	2478910	20.0
Naphthalene, 2,...	13.728	4.1	ng/u1	697713	3	14.551	3398670	20.0
Naphthalene, 2,...	13.869	5.7	ng/u1	965680	3	14.551	3398670	20.0
Naphthalene, 1,...	14.104	2.3	ng/u1	383968	3	14.551	3398670	20.0
2-Naphthaleneca...	14.675	3.4	ng/u1	572792	3	14.551	3398670	20.0
[1,1'-Biphenyl]...	16.034	3.1	ng/u1	602103	4	17.298	3855460	20.0
1-Acenaphthenol	16.263	24.5	ng/u1	4719120	4	17.298	3855460	20.0
Dibenzothiophene	17.116	2.3	ng/u1	439990	4	17.298	3855460	20.0
3-Nitrocarbazole	19.780	3.0	ng/u1	657426	5	21.480	4331070	20.0