LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
Data File : BM037748.D
Acq On : 27 Nov 2022 00:13
Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BHB42

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_M\Methods\SFAM-EPA-BM112322.M
Title : SVOA CALIBRATION

Signal : TIC: BM037748.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.469	11	14	26	rVB	101912	146091	0.93%	0.095%
2	3.875	80	83	97	rBV	328538	600540	3.81%	0.390%
3	4.110	118	123	130	rVB	34666	58688	0.37%	0.038%
4	6.087	454	459	465	rBV	48249	75383	0.48%	0.049%
5	6.575	536	542	548	rBV	67128	100780	0.64%	0.066%
6	7.257	652	658	665	rBV	392336	612560	3.89%	0.398%
7	7.428	679	687	694	rBV	1074751	1634968	10.38%	1.063%
8	7.634	711	722	729	rBV	1102887	1707503	10.84%	1.110%
9	8.110	795	803	809	rBV	926532	1462372	9.29%	0.951%
10	8.486	860	867	874	rVB	137951	223643	1.42%	0.145%
11	8.651	889	895	903	rVV	571730	910801	5.78%	0.592%
12	8.816	912	923	933	rBV	1003068	1671178	10.61%	1.087%
13	9.281	994	1002	1012	rVV	1404770	2418771	15.36%	1.573%
14	9.386	1014	1020	1026	rBV5	24159	56122	0.36%	0.036%
15	9.475	1026	1035	1042	rVV	1707602	2631072	16.71%	1.711%
16	9.598	1049	1056	1062	rVV	408333	907684	5.76%	0.590%
17	9.663	1062	1067	1078	rVV	639006	1050875	6.67%	0.683%
18	10.004	1118	1125	1135	rBV	1135919	1800601	11.44%	1.171%
19	10.198	1155	1158	1168	rVB3	35336	68689	0.44%	0.045%
20	10.328	1173	1180	1191	rVB2	91384	173365	1.10%	0.113%
21	10.545	1210	1217	1227	rVB	1603100	2637171	16.75%	1.715%
22	10.839	1262	1267	1271	rVV	46554	80575	0.51%	0.052%
23	10.927	1275	1282	1288	rVV	1357530	2322026	14.75%	1.510%
24	10.986	1288	1292	1298	rVV3	120558	214318	1.36%	0.139%
25	11.063	1298	1305	1307	rVV	1110948	1722192	10.94%	1.120%
26	11.104	1307	1312	1320	rVV	5002108	8534025	54.20%	5.548%
27	11.169	1320	1323	1328	rVV	74703	126164	0.80%	0.082%
28	11.222	1328	1332	1337	rVV6	32861	68423	0.43%	0.044%
29	11.275	1337	1341	1348	rVV6	38475	98673	0.63%	0.064%
30	11.339	1348	1352	1359	rVV3	37022	82090	0.52%	0.053%
31	11.510	1372	1381	1389	rVB9	28940	98882	0.63%	0.064%
32	11.663	1401	1407	1410	rBV	46097	74358	0.47%	0.048%
33	11.698	1410	1413	1419	rVV3	34286	55449	0.35%	0.036%
34	11.839	1431	1437	1443	rVB	49477	80793	0.51%	0.053%
35	11.963	1447	1458	1464	rBV	61874	130335	0.83%	0.085%
36	12.039	1464	1471	1482	rVB	575582	1188591	7.55%	0.773%

LSC Area Percent Report

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Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BHB42

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_M\Methods\SFAM-EPA-BM112322.M
Title : SVOA CALIBRATION

37	12.427	1531	1537	1540	rBV5	33777	64645	0.41%	0.042%
38	12.474	1540	1545	1549	rBV	321691	522440	3.32%	0.340%
39	12.792	1590	1599	1604	rBV2	113310	276223	1.75%	0.180%
40	12.886	1610	1615	1623	rVB	121068	177479	1.13%	0.115%
41	13.080	1643	1648	1655	rVB4	30991	56609	0.36%	0.037%
42	13.551	1720	1728	1735	rBV4	47688	131424	0.83%	0.085%
43	13.616	1735	1739	1749	rVB5	47336	99031	0.63%	0.064%
44	13.698	1749	1753	1762	rBV9	26864	66513	0.42%	0.043%
45	14.139	1822	1828	1834	rBV	3165068	4144548	26.32%	2.695%
46	14.433	1869	1878	1888	rVV	3427397	5586349	35.48%	3.632%
47	14.733	1921	1929	1935	rBV	2250708	3192809	20.28%	2.076%
48	14.792	1935	1939	1946	rVV4	47942	88074	0.56%	0.057%
49	14.921	1955	1961	1973	rVB	465608	706912	4.49%	0.460%
50	15.127	1991	1996	2002	rVV	231758	329754	2.09%	0.214%
51	15.210	2005	2010	2019	rVV	240793	406689	2.58%	0.264%
52	15.292	2019	2024	2028	rBV	104619	170114	1.08%	0.111%
53	15.333	2028	2031	2038	rVB	117172	195174	1.24%	0.127%
54	15.486	2052	2057	2064	rVB9	46686	113743	0.72%	0.074%
55	15.586	2067	2074	2077	rBV	112391	196142	1.25%	0.128%
56	15.721	2089	2097	2103	rVV	4540155	6439984	40.90%	4.187%
57	15.792	2103	2109	2112	rVV	1629621	2481444	15.76%	1.613%
58	15.827	2112	2115	2120	rVV	1652068	2139700	13.59%	1.391%
59	15.862	2120	2121	2130	rVB3	124806	218073	1.38%	0.142%
60	16.080	2151	2158	2164	rVB2	273031	459215	2.92%	0.299%
61	16.168	2168	2173	2177	rBV	134454	194618	1.24%	0.127%
62	16.209	2177	2180	2184	rVB3	72072	92153	0.59%	0.060%
63	16.351	2200	2204	2208	rVB2	78786	107227	0.68%	0.070%
64	16.468	2209	2224	2233	rBV	3834633	9470123	60.14%	6.157%
65	16.651	2248	2255	2259	rBV	3217793	4986174	31.67%	3.242%
66	16.692	2259	2262	2265	rVV2	789284	1117800	7.10%	0.727%
67	16.721	2265	2267	2270	rVB2	209422	233033	1.48%	0.152%
68	16.874	2289	2293	2300	rVB3	72019	119377	0.76%	0.078%
69	17.051	2309	2323	2337	rBV2	2306302	6528840	41.46%	4.245%
70	17.186	2340	2346	2350	rVV	212597	411932	2.62%	0.268%
71	17.233	2350	2354	2363	rVB3	305669	598152	3.80%	0.389%
72	17.380	2369	2379	2389	rBV2	5236260	15746364	100.00%	10.238%
73	17.468	2389	2394	2402	rVV	2986938	4754720	30.20%	3.091%
74	17.568	2402	2411	2428	rVV	5823619	9644318	61.25%	6.270%
75	17.692	2428	2432	2435	rVV	407599	610867	3.88%	0.397%
76	17.727	2435	2438	2445	rVB	310604	486455	3.09%	0.316%

LSC Area Percent Report

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Data File : BM037748.D
Acq On : 27 Nov 2022 00:13
Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHB42

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

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Title : SVOA CALIBRATION

77	17.839	2449	2457	2465	rBV4	869354	2274004	14.44%	1.478%
78	17.909	2465	2469	2475	rVB2	274408	534476	3.39%	0.347%
79	17.998	2480	2484	2491	rVB2	216567	356623	2.26%	0.232%
80	18.092	2496	2500	2503	rBV	97869	123884	0.79%	0.081%
81	18.133	2503	2507	2512	rVB	233732	294410	1.87%	0.191%
82	18.180	2512	2515	2518	rVB	65802	79962	0.51%	0.052%
83	18.245	2519	2526	2530	rBV2	274212	581001	3.69%	0.378%
84	18.356	2540	2545	2553	rBV3	561534	923615	5.87%	0.600%
85	18.427	2553	2557	2571	rVB8	108435	333137	2.12%	0.217%
86	18.609	2584	2588	2592	rBV	121309	181184	1.15%	0.118%
87	18.692	2599	2602	2612	rVV2	142146	270253	1.72%	0.176%
88	18.768	2612	2615	2620	rVB2	156161	180841	1.15%	0.118%
89	19.245	2690	2696	2705	rBV2	648330	1082306	6.87%	0.704%
90	19.486	2733	2737	2741	rVB	66246	101135	0.64%	0.066%
91	19.615	2755	2759	2769	rVB	274684	407806	2.59%	0.265%
92	19.850	2793	2799	2805	rBV	6811001	8831895	56.09%	5.742%
93	20.250	2864	2867	2872	rVB	131876	151558	0.96%	0.099%
94	20.774	2953	2956	2958	rBV2	86774	94570	0.60%	0.061%
95	20.821	2962	2964	2967	rVB	97041	93660	0.59%	0.061%
96	21.321	3045	3049	3056	rBV	613464	786116	4.99%	0.511%
97	21.633	3097	3102	3109	rBV	3597141	4421140	28.08%	2.874%
98	22.909	3315	3319	3326	rVB	437971	616160	3.91%	0.401%
99	23.962	3491	3498	3512	rVB2	3823221	8038750	51.05%	5.226%
100	24.127	3519	3526	3534	rVB2	1899620	3859645	24.51%	2.509%

Sum of corrected areas: 153809123

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
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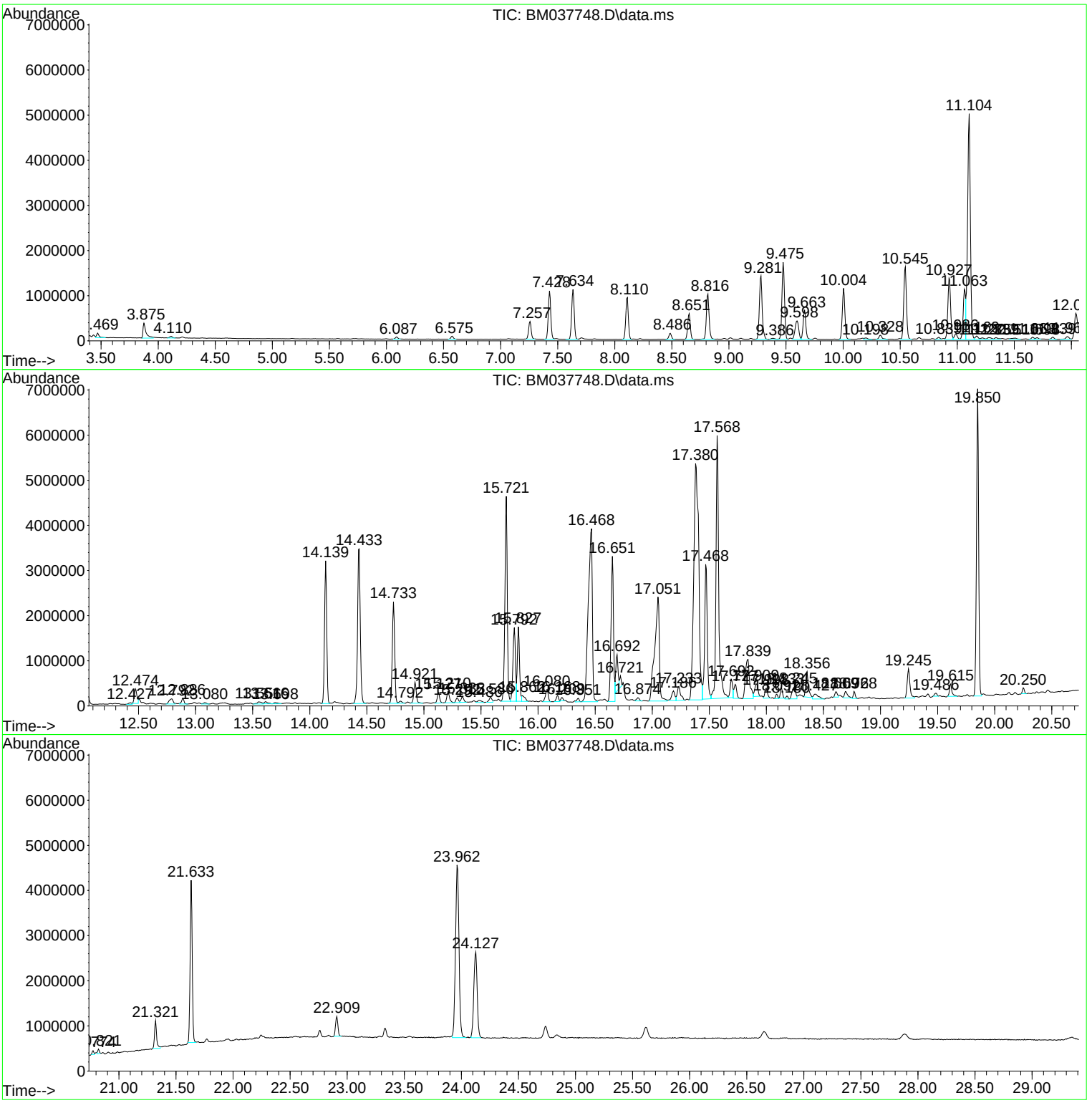
Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHB42

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
Data File : BM037748.D
Acq On : 27 Nov 2022 00:13
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Misc :
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Instrument :
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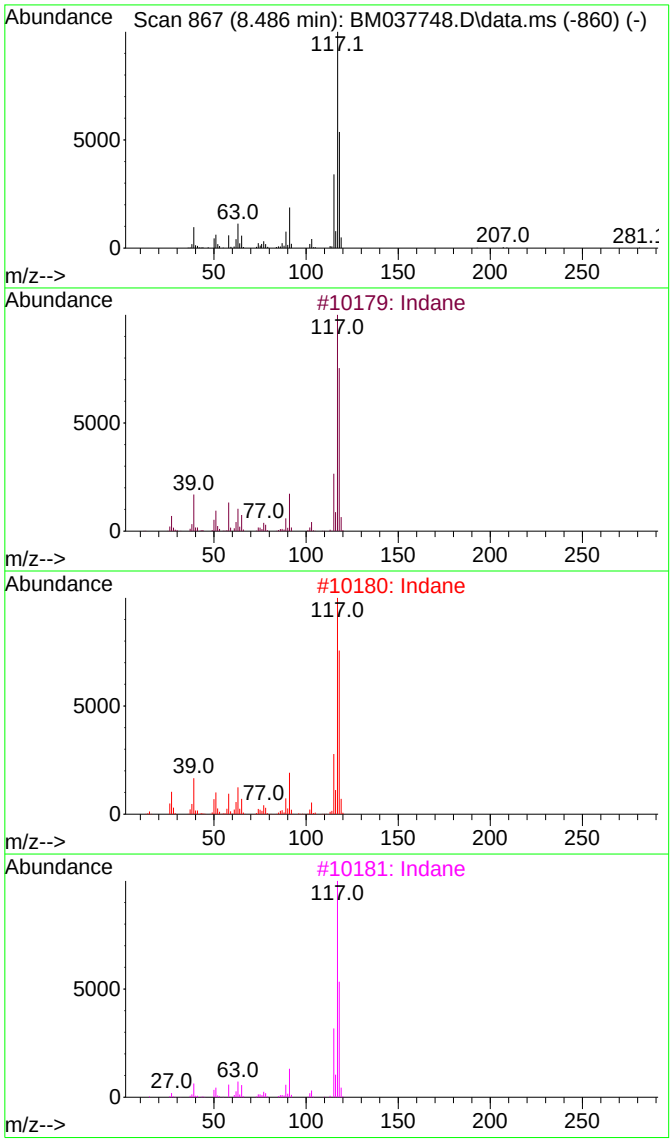
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

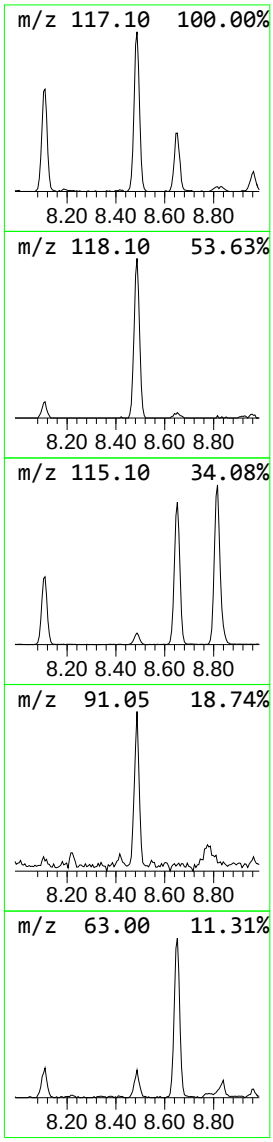
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Indane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.486	3.06 ng/ul	223643	1,4-Dichlorobenzene-d4	8.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	90
2	Indane	118	C9H10	000496-11-7	90
3	Indane	118	C9H10	000496-11-7	87
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	78
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	74





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
Data File : BM037748.D
Acq On : 27 Nov 2022 00:13
Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHB42

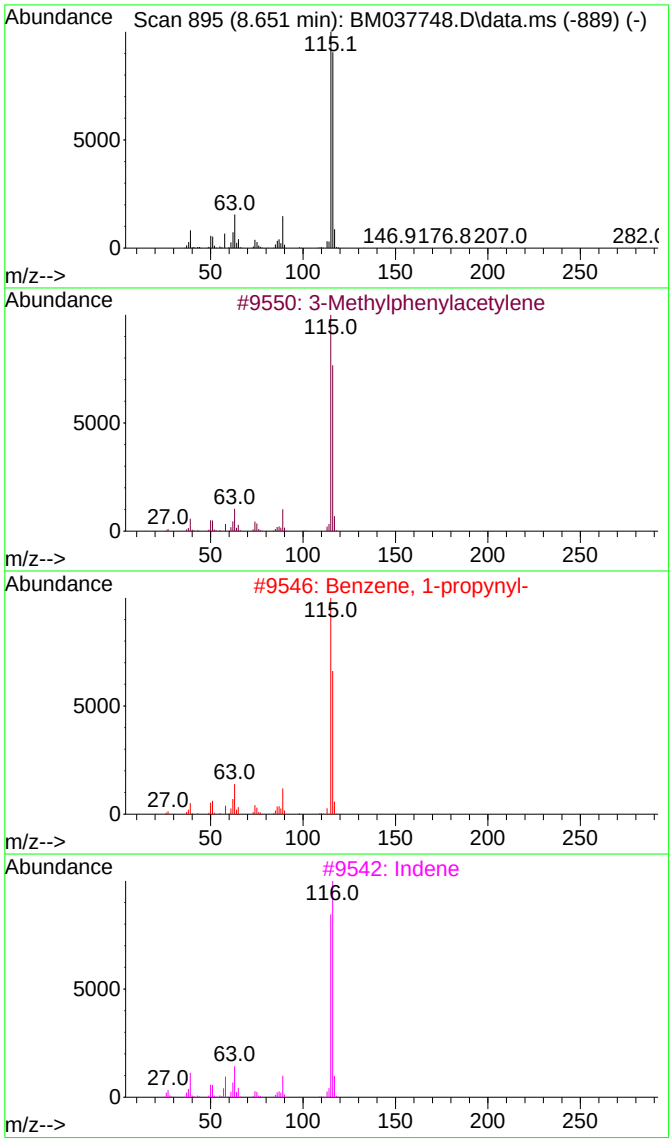
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

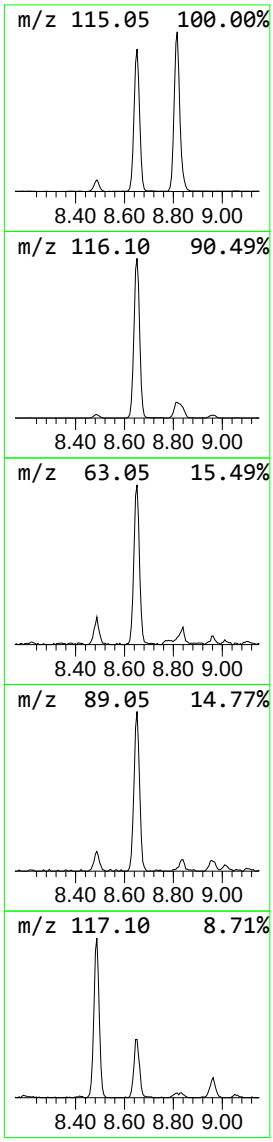
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 3-Methylphenylacetylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.651	12.46 ng/ul	910801	1,4-Dichlorobenzene-d4	8.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methylphenylacetylene	116	C9H8	000766-82-5	94
2	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3	Indene	116	C9H8	000095-13-6	94
4	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
Data File : BM037748.D
Acq On : 27 Nov 2022 00:13
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ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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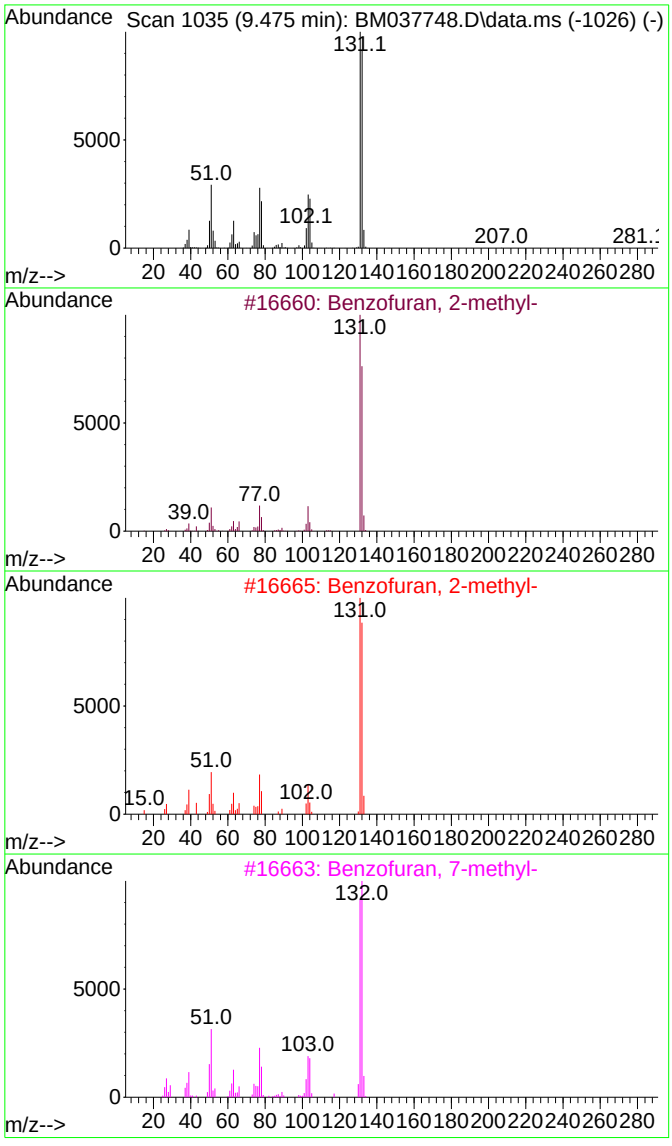
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

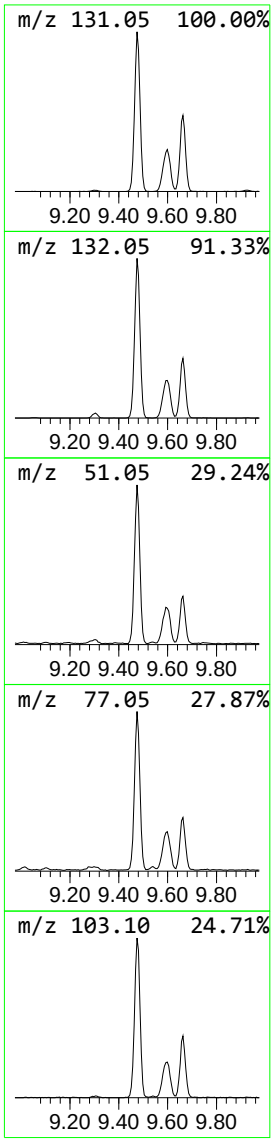
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzofuran, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.475	35.98 ng/ul	2631070	1,4-Dichlorobenzene-d4	8.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	95
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
4	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	93
5	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	93





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ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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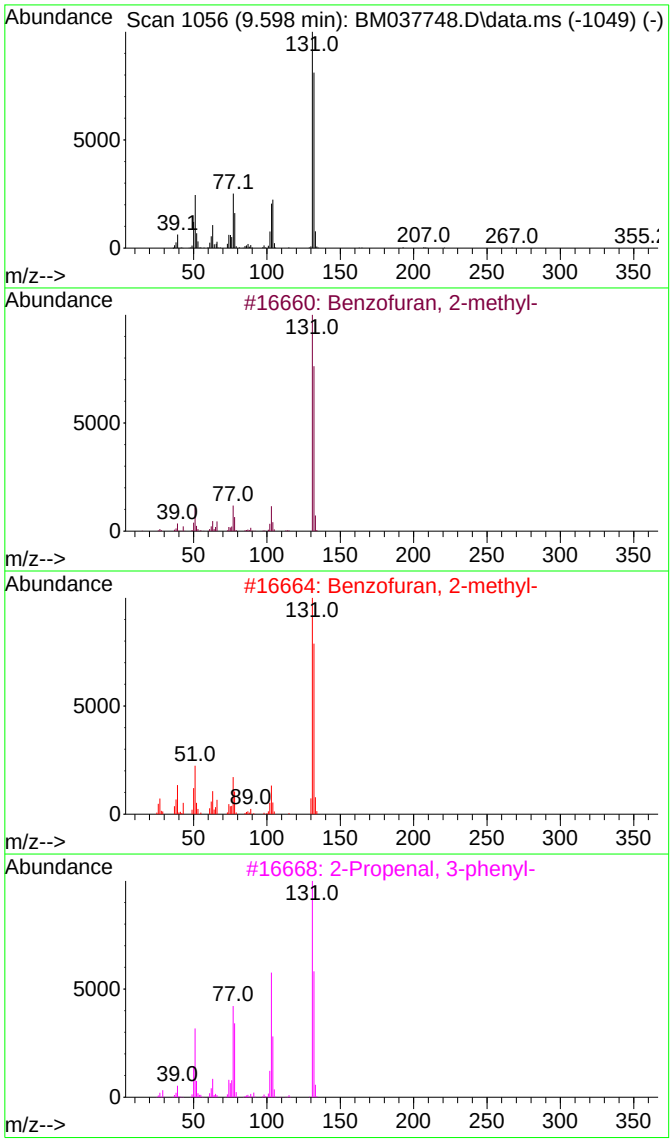
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

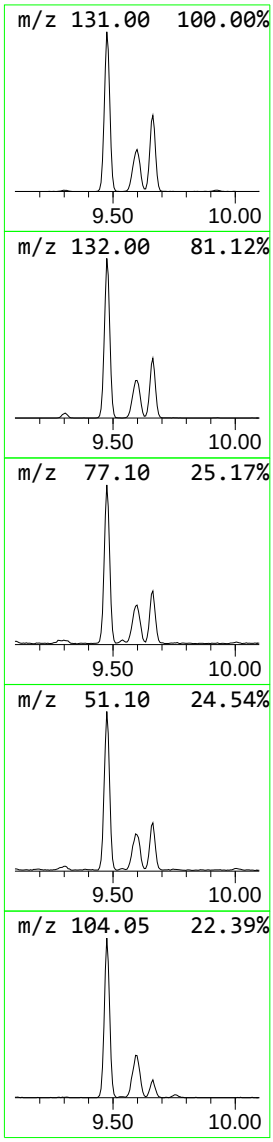
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 2-Propenal, 3-phenyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.598	7.82 ng/ul	907684	Naphthalene-d8	10.928

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	93
4	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91
5	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
Data File : BM037748.D
Acq On : 27 Nov 2022 00:13
Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHB42

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

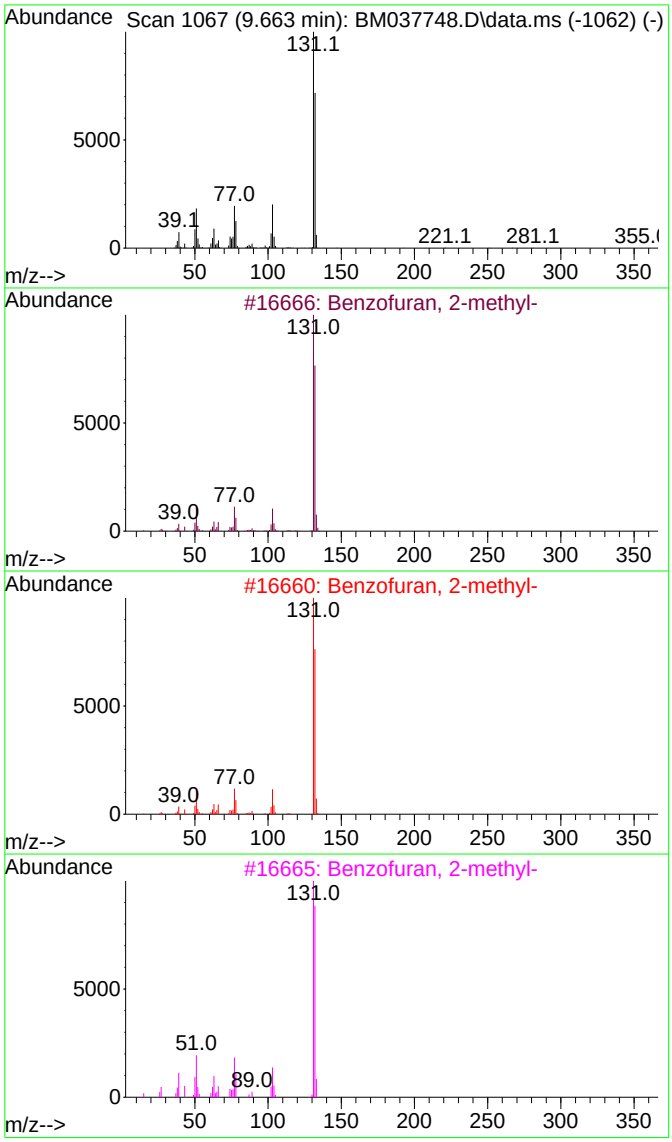
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

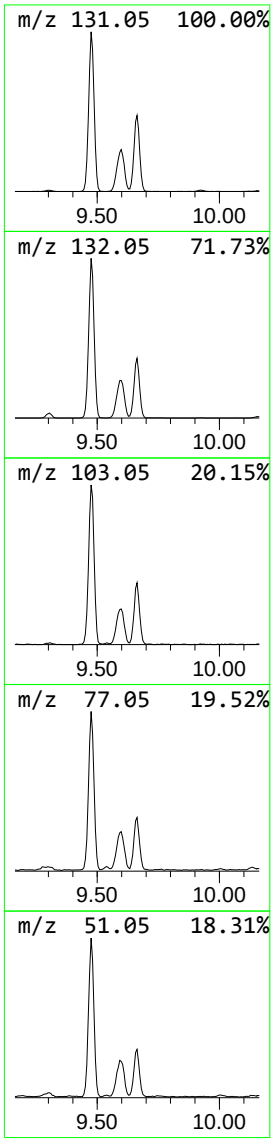
****

Peak Number 5 3-Phenyl-2-propyn-1-ol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.663	9.05 ng/ul	1050880	Naphthalene-d8	10.928

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
5	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	87





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Acq On : 27 Nov 2022 00:13
Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHB42

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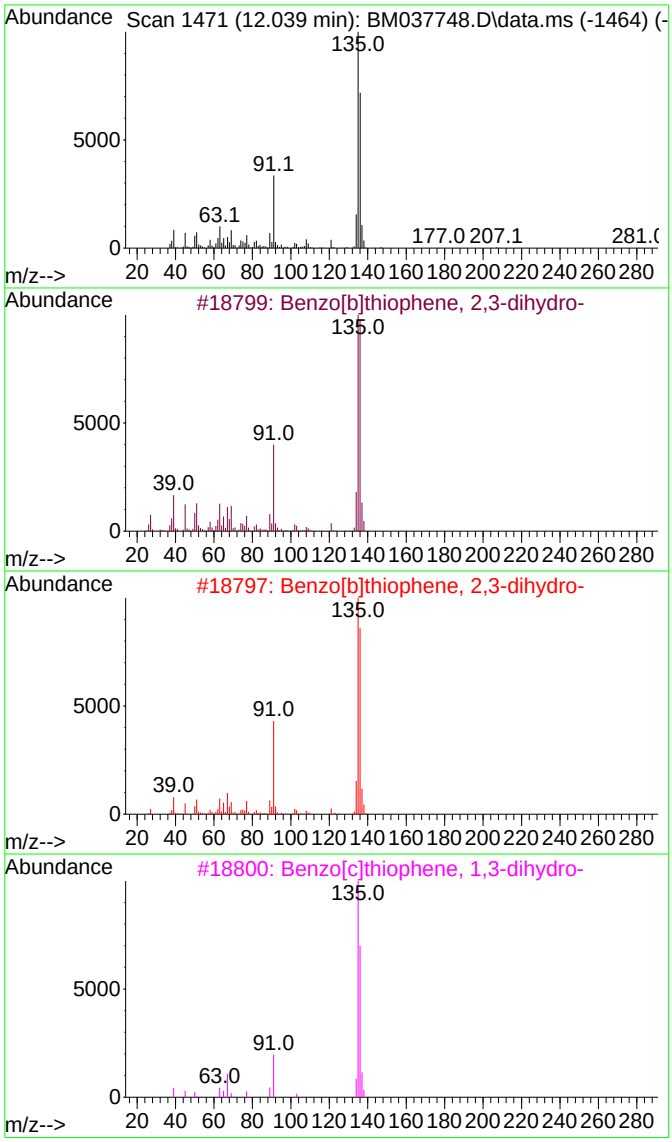
Quant Title : SVOA CALIBRATION

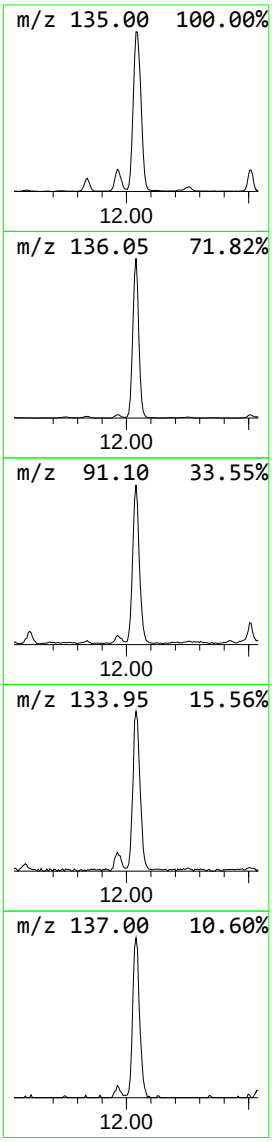
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzo[b]thiophene, 2,3-dihy... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.039	10.24 ng/ul	1188590	Naphthalene-d8	10.928

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	94
2	Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	93
3	Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	86
4	Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	81
5	Benzene, (ethenylthio)-	136	C8H8S	001822-73-7	74





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Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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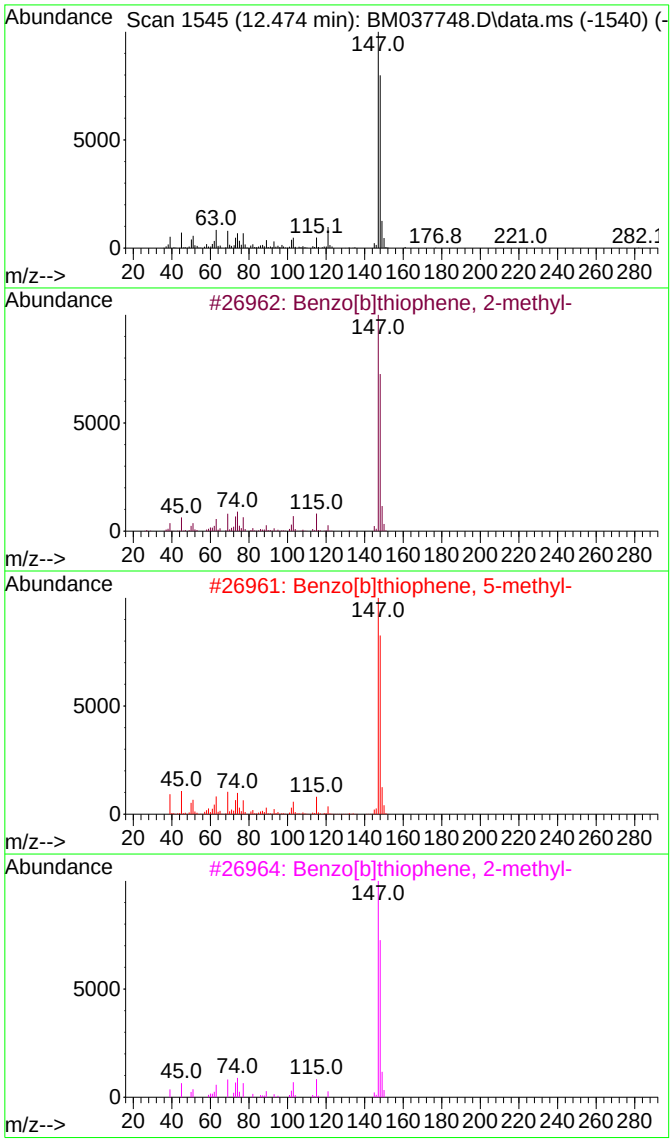
Quant Title : SVOA CALIBRATION

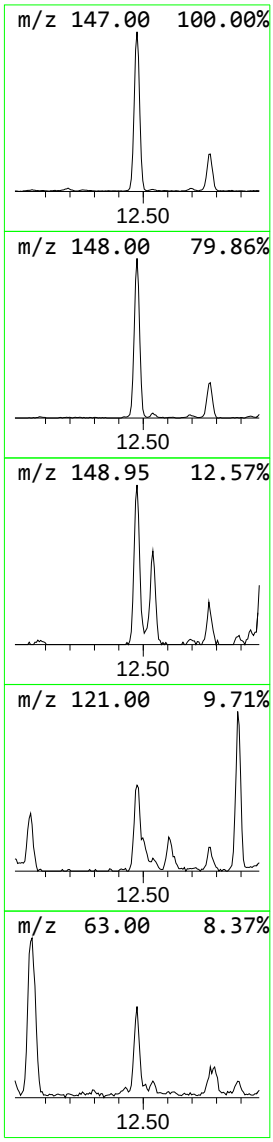
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzo[b]thiophene, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.474	4.50 ng/ul	522440	Naphthalene-d8	10.928

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
2	Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	90
3	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	90
4	3-Methylbenzothiophene	148	C9H8S	001455-18-1	90
5	Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	90





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ClientSampleId :
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Quant Title : SVOA CALIBRATION

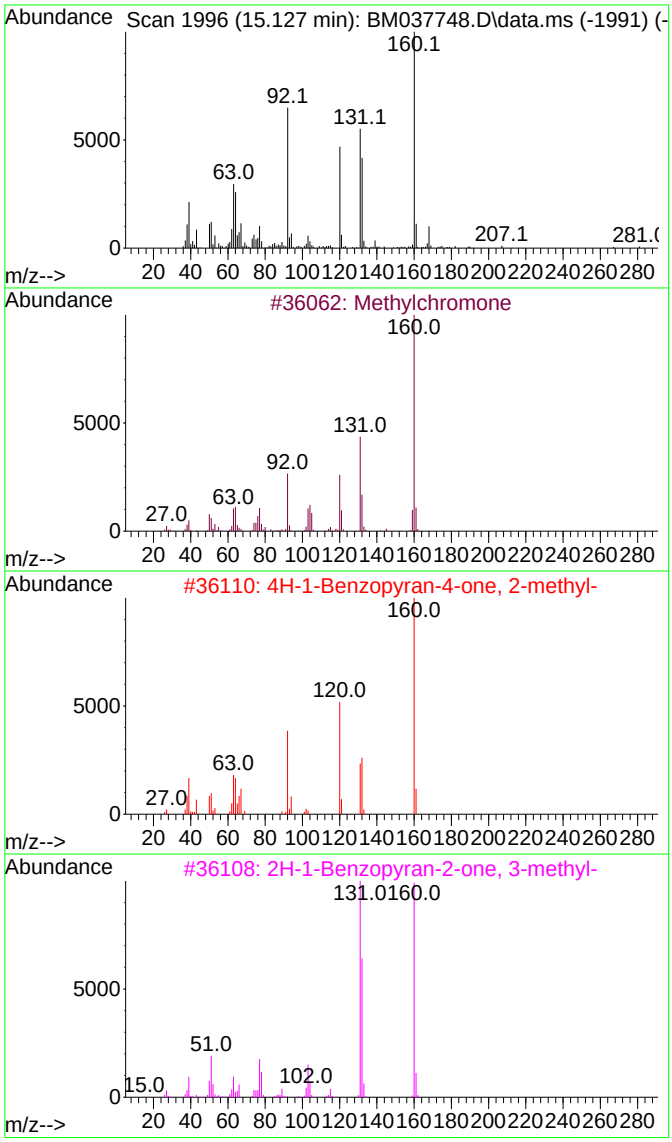
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

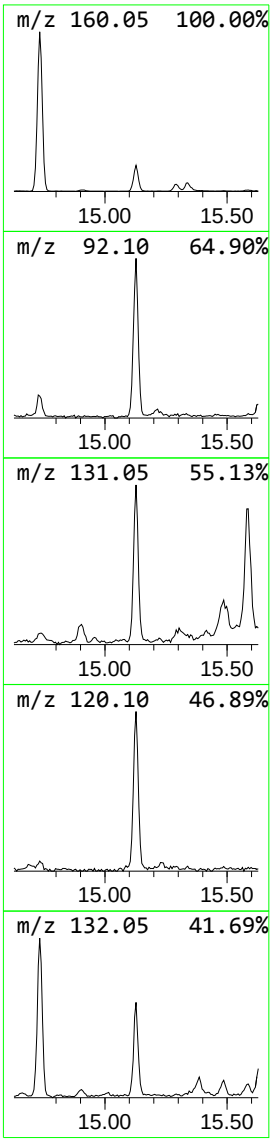
****

Peak Number 9 Methylchromone Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.127	2.07 ng/ul	329754	Acenaphthene-d10	14.733

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methylchromone	160	C10H8O2	000085-90-5	72
2	4H-1-Benzopyran-4-one, 2-methyl-	160	C10H8O2	005751-48-4	64
3	2H-1-Benzopyran-2-one, 3-methyl-	160	C10H8O2	002445-82-1	55
4	2H-1-Benzopyran-2-one, 3-methyl-	160	C10H8O2	002445-82-1	49
5	2H-1-Benzopyran-2-one, 4-methyl-	160	C10H8O2	000607-71-6	49





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Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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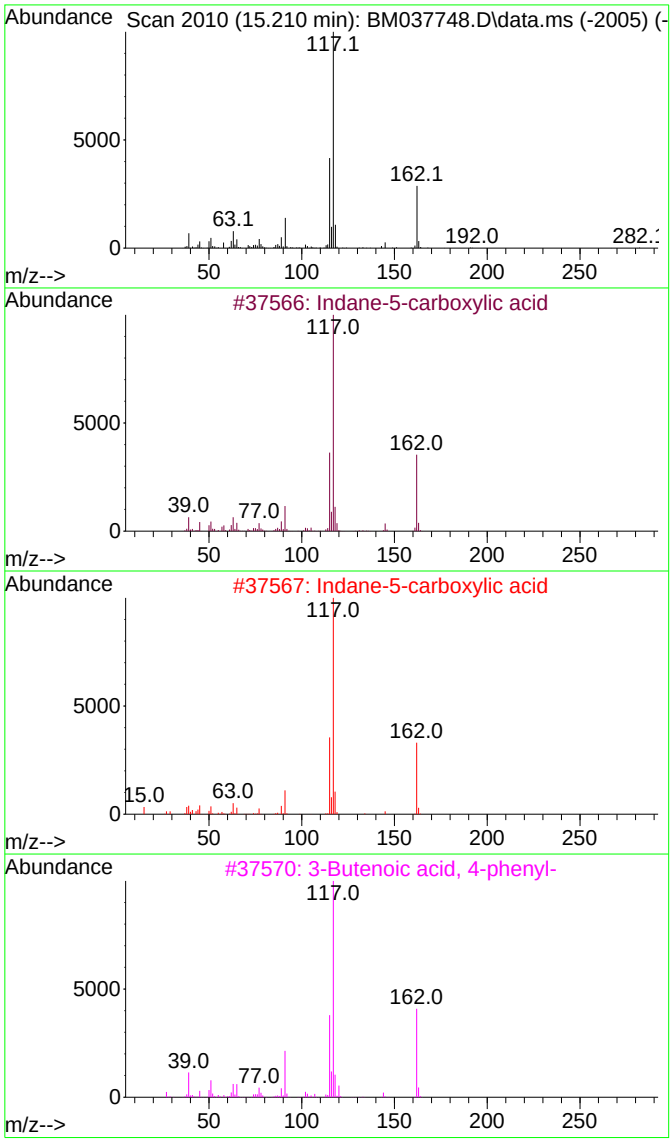
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

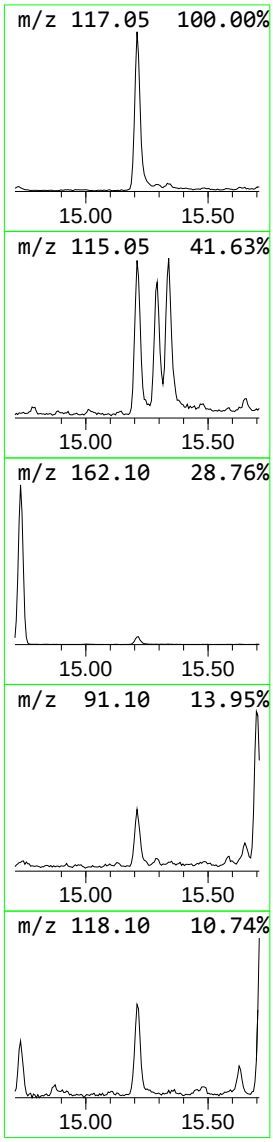
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Indane-5-carboxylic acid Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.210	2.55 ng/ul	406689	Acenaphthene-d10	14.733

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane-5-carboxylic acid	162	C10H10O2	065898-38-6	93
2	Indane-5-carboxylic acid	162	C10H10O2	065898-38-6	87
3	3-Butenoic acid, 4-phenyl-	162	C10H10O2	002243-53-0	80
4	trans-4-Phenyl-3-butenoic acid	162	C10H10O2	001914-58-5	76
5	Benzene, (3-chloroallyl)-	152	C9H9Cl	006268-37-7	64





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ClientSampleId :
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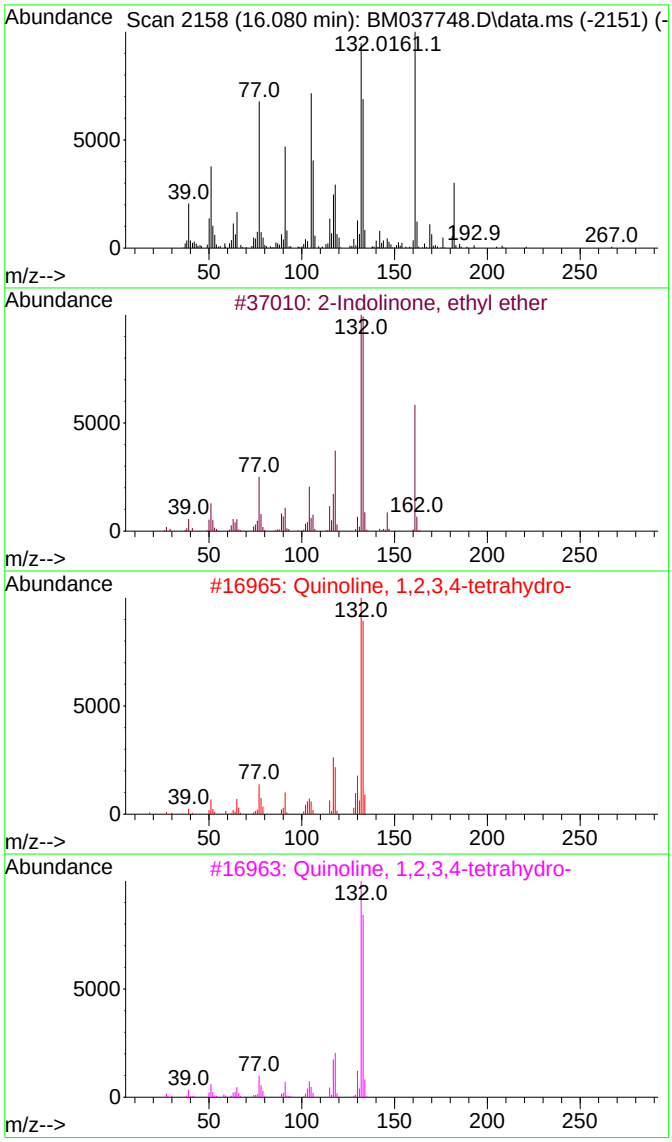
Quant Title : SVOA CALIBRATION

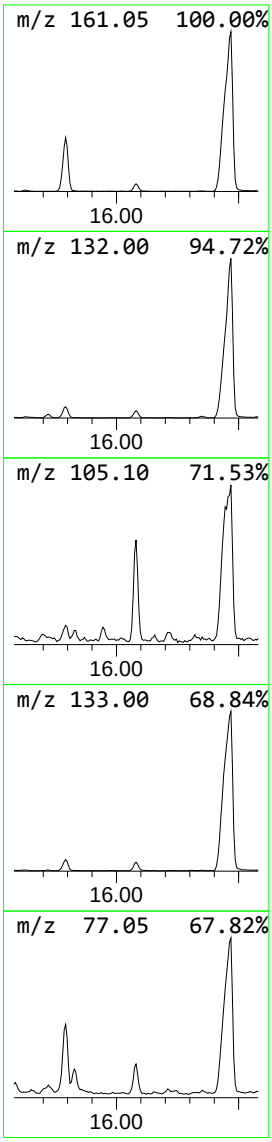
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 2-Indolinone, ethyl ether Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.080	2.88 ng/ul	459215	Acenaphthene-d10	14.733

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Indolinone, ethyl ether	161	C10H11NO	1000453-17-4	76
2	Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	55
3	Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	55
4	Quinoline, 5,6,7,8-tetrahydro-	133	C9H11N	010500-57-9	46
5	Isoquinoline, 5,6,7,8-tetrahydro-	133	C9H11N	036556-06-6	45





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Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

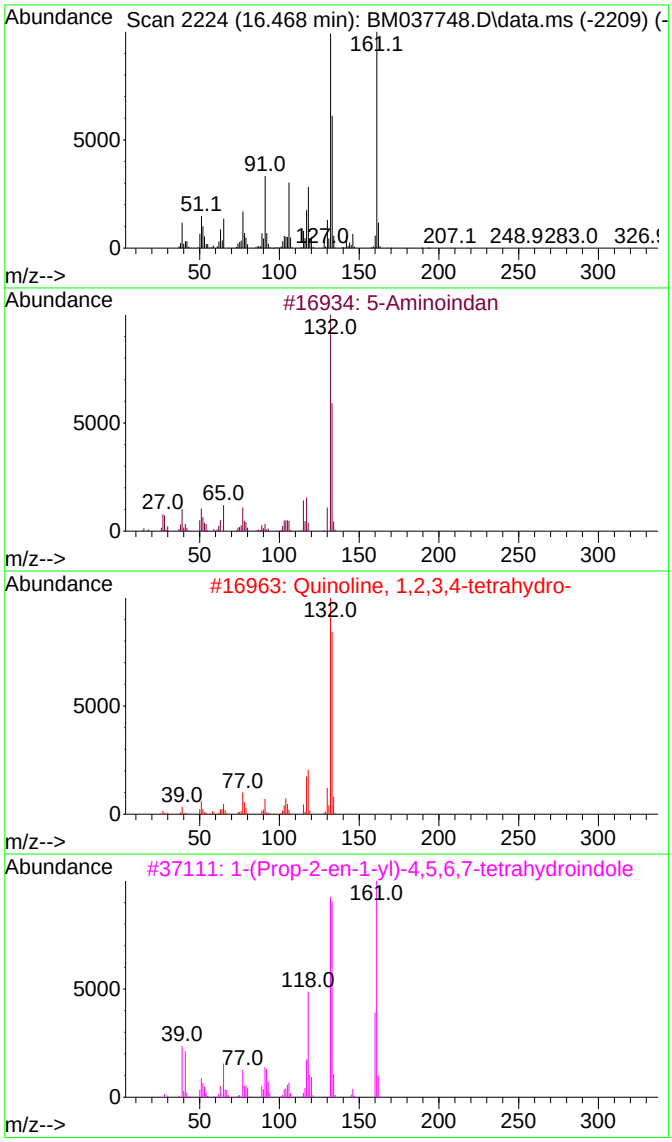
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

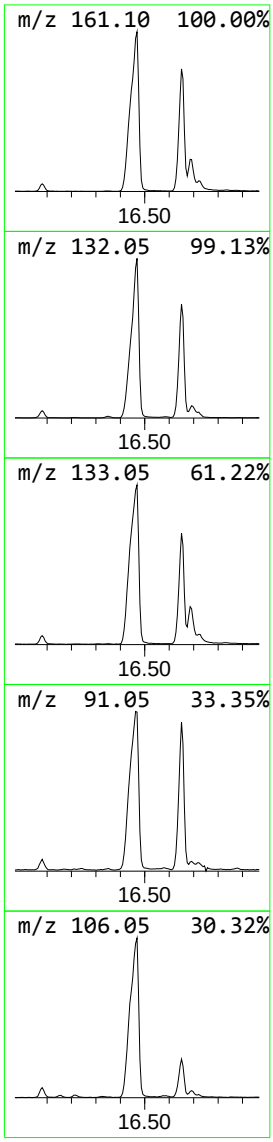
****

Peak Number 12 5-Aminoindan Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.468	39.83 ng/ul	9470120	Phenanthrene-d10	17.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Aminoindan	133	C9H11N	024425-40-9	86
2	Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	55
3	1-(Prop-2-en-1-yl)-4,5,6,7-tetra...	161	C11H15N	1450892-88-2	52
4	Quinoline, 5,6,7,8-tetrahydro-	133	C9H11N	010500-57-9	50
5	Benzenamine, 2-cyclopropyl-	133	C9H11N	1000327-33-3	46





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Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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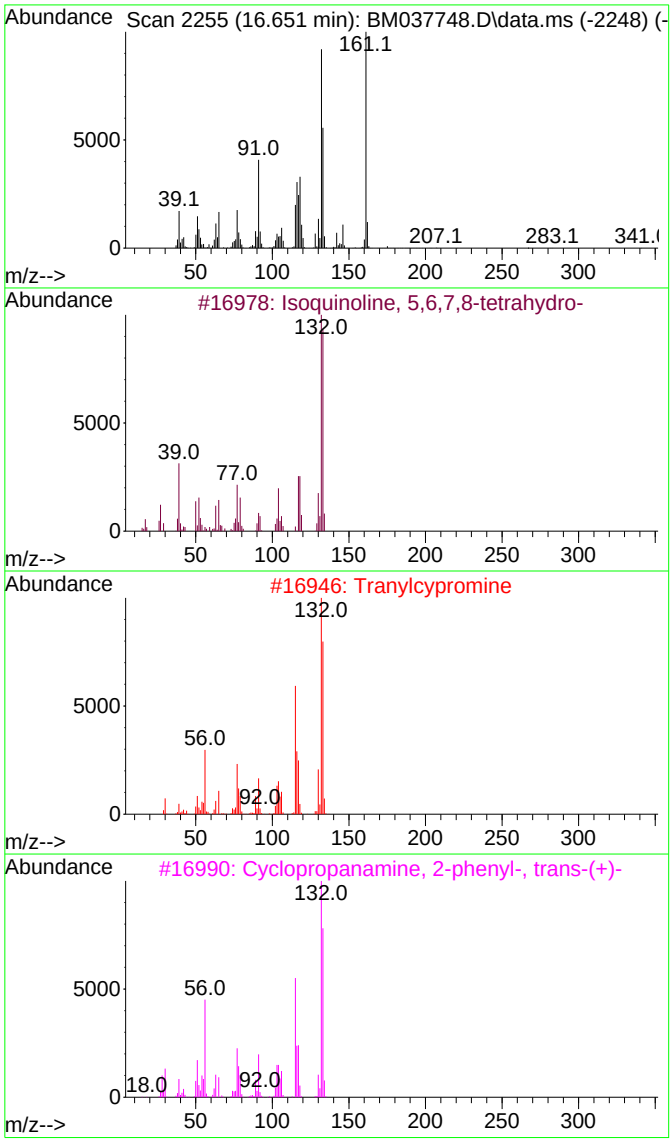
Quant Title : SVOA CALIBRATION

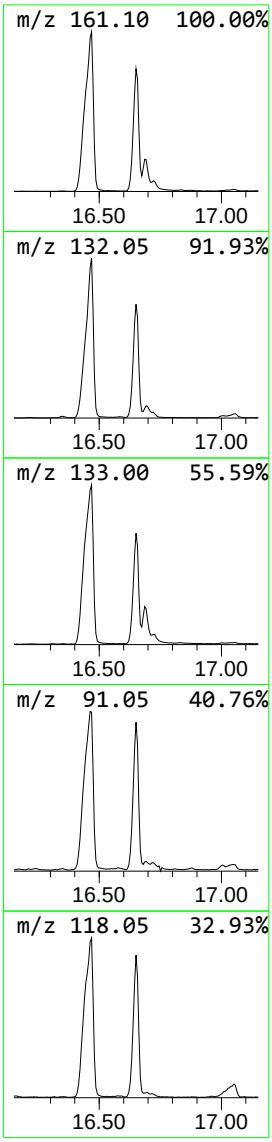
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Isoquinoline, 5,6,7,8-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.651	20.97 ng/ul	4986170	Phenanthrene-d10	17.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isoquinoline, 5,6,7,8-tetrahydro-	133	C9H11N	036556-06-6	78
2	Tranylcypromine	133	C9H11N	000155-09-9	55
3	Cyclopropanamine, 2-phenyl-, tra...	133	C9H11N	003721-28-6	55
4	Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	50
5	Tranylcypromine	133	C9H11N	000155-09-9	49





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Misc :
ALS Vial : 60 Sample Multiplier: 1

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ClientSampleId :
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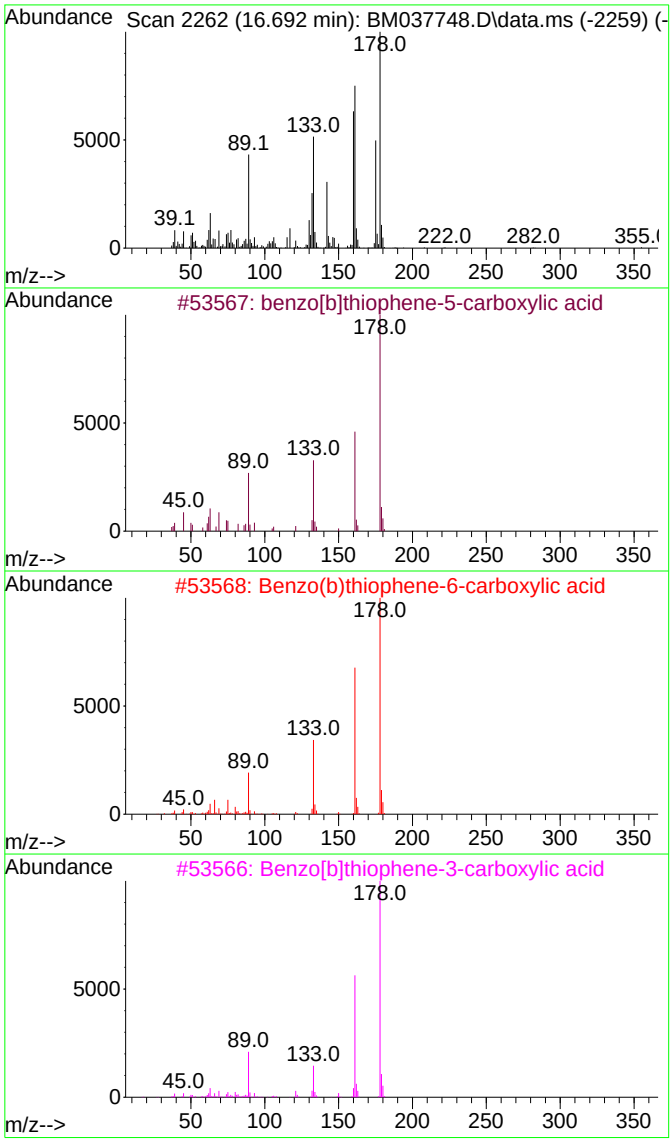
Quant Title : SVOA CALIBRATION

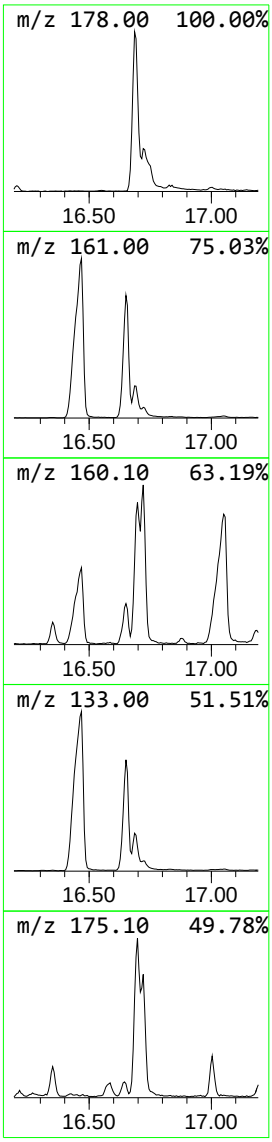
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 benzo[b]thiophene-5-carboxy... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.692	4.70 ng/ul	1117800	Phenanthrene-d10	17.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	benzo[b]thiophene-5-carboxylic acid	178	C9H6O2S	1000404-77-7	90
2	Benzo(b)thiophene-6-carboxylic acid	178	C9H6O2S	006179-26-6	70
3	Benzo[b]thiophene-3-carboxylic acid	178	C9H6O2S	005381-25-9	46
4	2,3,4,5-Tetramethylbenzoic acid	178	C11H14O2	002529-39-7	43
5	3-Methyl-7,8-dihydroquinolin-5(6...	161	C10H11NO	060247-70-3	42





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ClientSampleId :
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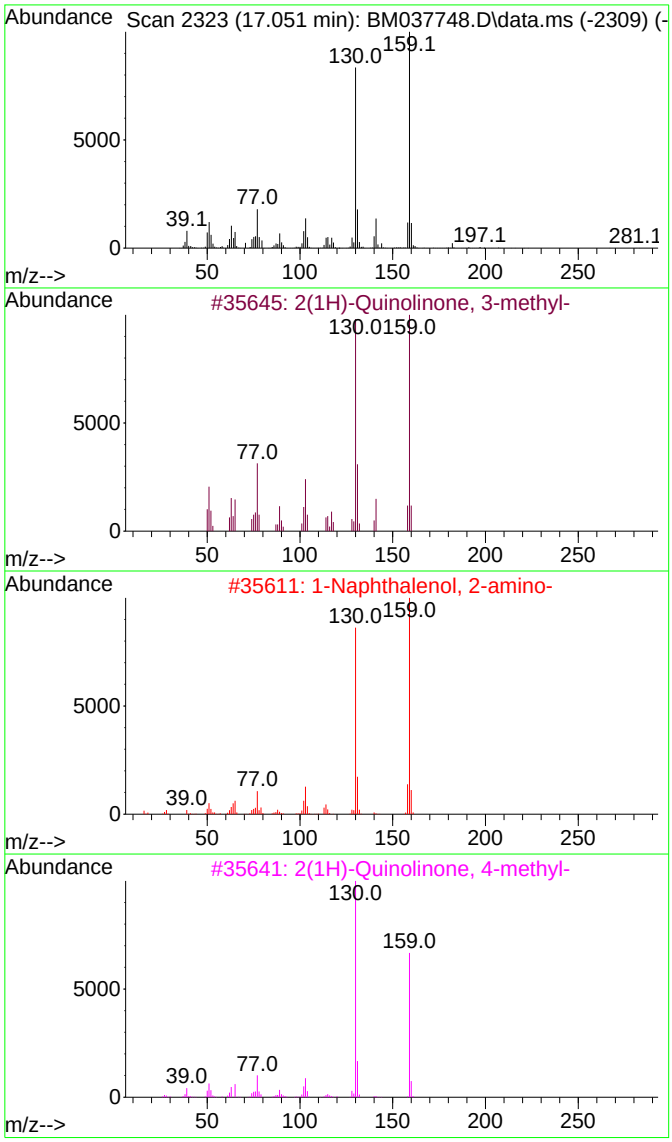
Quant Title : SVOA CALIBRATION

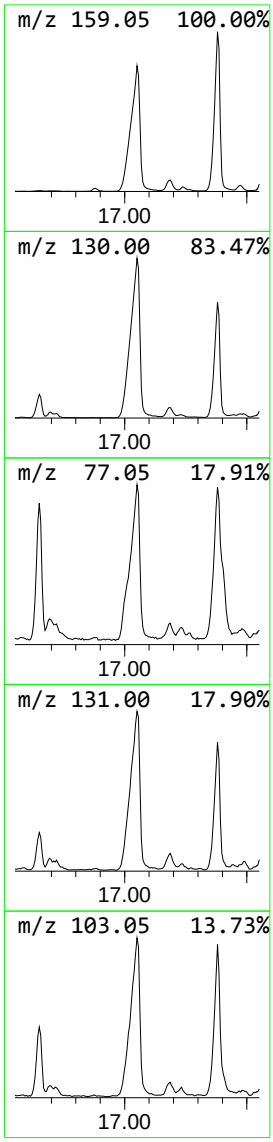
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 2(1H)-Quinolinone, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.051	27.46 ng/ul	6528840	Phenanthrene-d10	17.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Quinolinone, 3-methyl-	159	C10H9NO	002721-59-7	97
2	1-Naphthalenol, 2-amino-	159	C10H9NO	000606-41-7	94
3	2(1H)-Quinolinone, 4-methyl-	159	C10H9NO	000607-66-9	91
4	2(1H)-Quinolinone, 4-methyl-	159	C10H9NO	000607-66-9	91
5	2-(Aminomethylidene)-2,3-dihydro...	159	C10H9NO	053394-92-6	90





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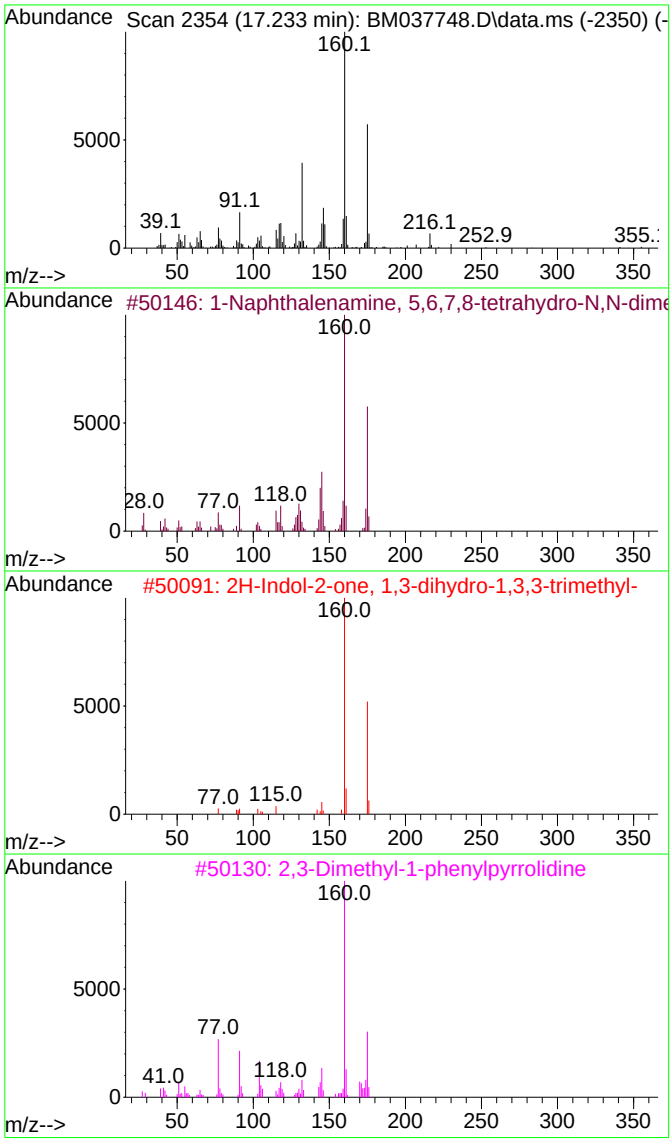
Quant Title : SVOA CALIBRATION

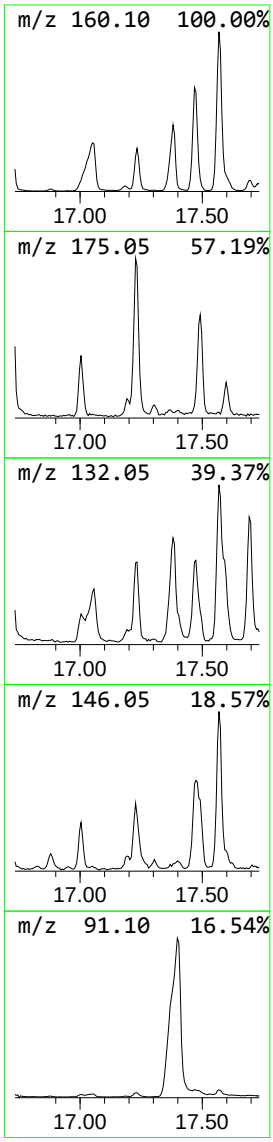
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 1-Naphthalenamine, 5,6,7,8-... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.233	2.52 ng/ul	598152	Phenanthrene-d10	17.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenamine, 5,6,7,8-tetra...	175	C12H17N	013541-35-0	58
2	2H-Indol-2-one, 1,3-dihydro-1,3,...	175	C11H13NO	020200-86-6	53
3	2,3-Dimethyl-1-phenylpyrrolidine	175	C12H17N	003783-58-2	50
4	1H-Isoindole, 3-methoxy-4,7-dime...	175	C11H13NO	100813-60-3	49
5	4-Hydroxy-7-methyl-1,8-naphthyri...	204	C10H8N2O3	013250-97-0	47





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Quant Title : SVOA CALIBRATION

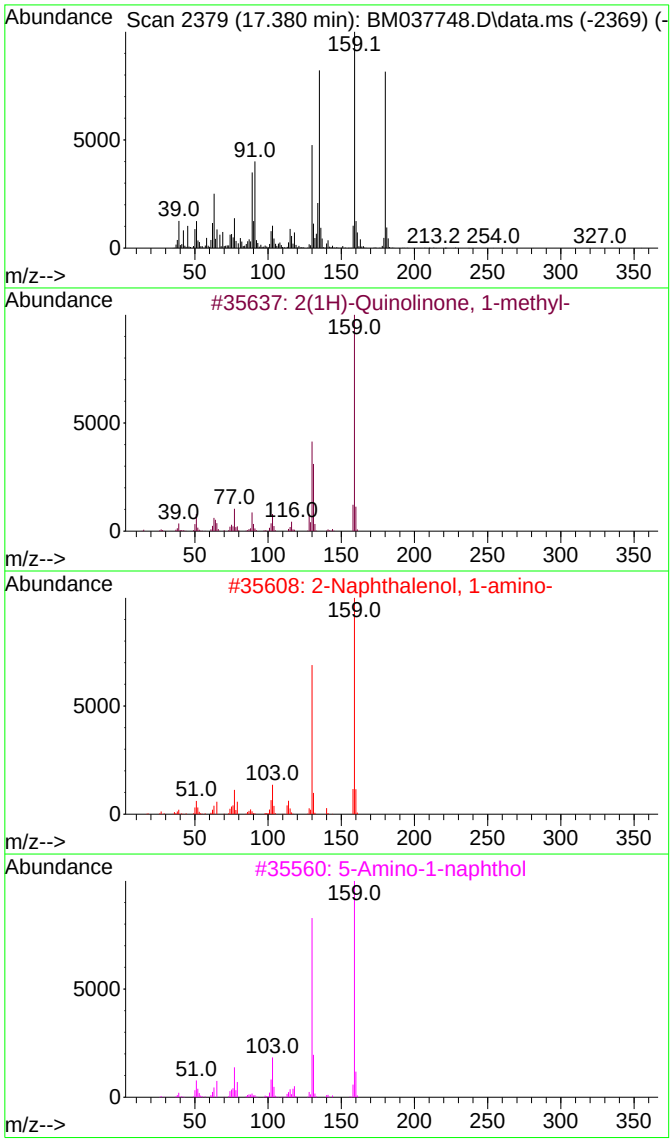
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

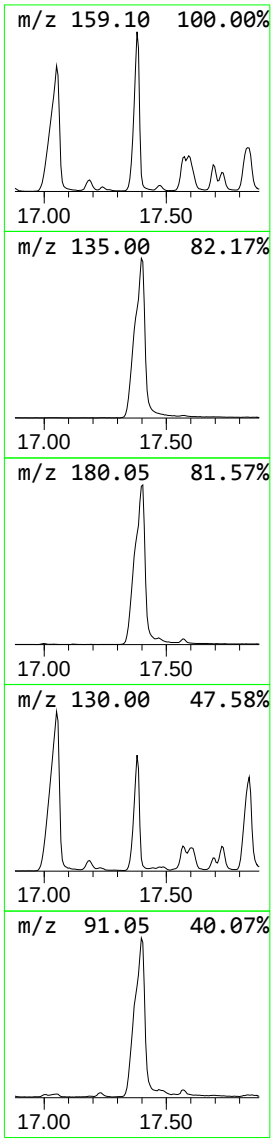
****

Peak Number 17 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.380	66.23 ng/ul	15746400	Phenanthrene-d10	17.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Quinolinone, 1-methyl-	159	C10H9NO	000606-43-9	38
2	2-Naphthalenol, 1-amino-	159	C10H9NO	002834-92-6	38
3	5-Amino-1-naphthol	159	C10H9NO	000083-55-6	38
4	2(1H)-Quinolinone, 3-methyl-	159	C10H9NO	002721-59-7	38
5	4-Quinolinol, 2-methyl-	159	C10H9NO	000607-67-0	38





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Data File : BM037748.D
Acq On : 27 Nov 2022 00:13
Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHB42

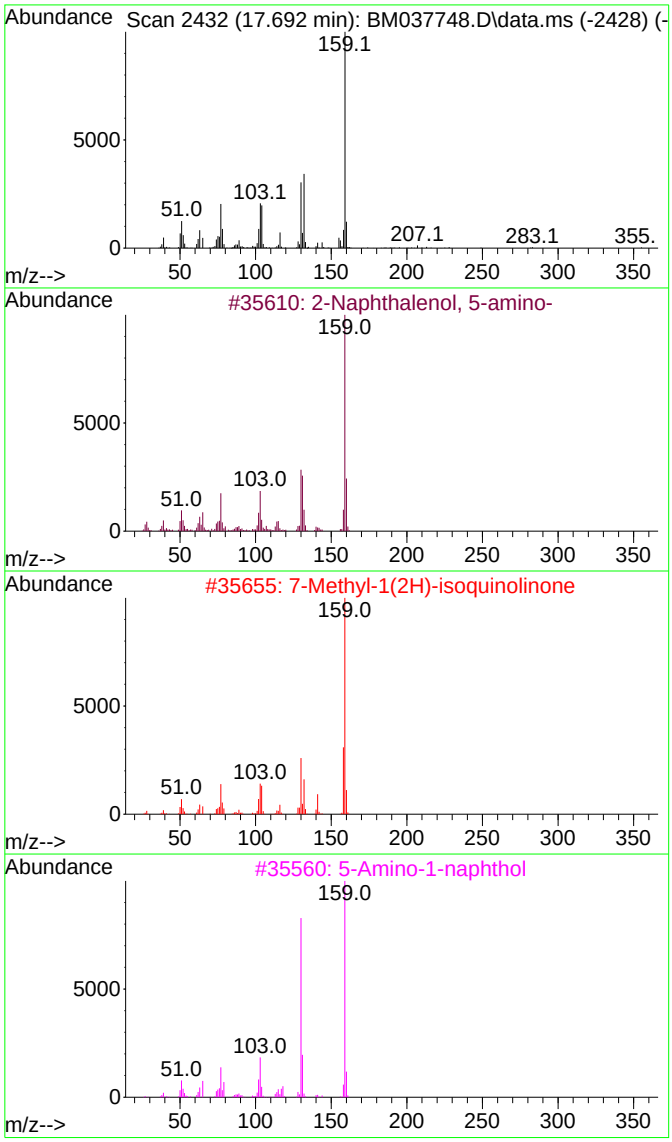
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

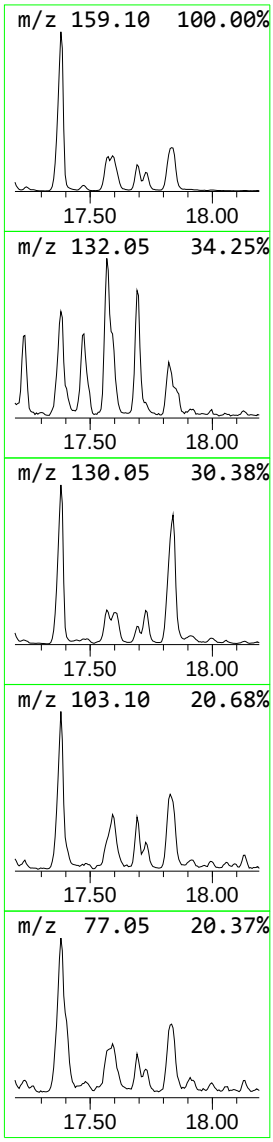
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 2-Naphthalenol, 5-amino- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.692	2.57 ng/ul	610867	Phenanthrene-d10	17.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenol, 5-amino-	159	C10H9NO	000086-97-5	76
2	7-Methyl-1(2H)-isoquinolinone	159	C10H9NO	026829-47-0	74
3	5-Amino-1-naphthol	159	C10H9NO	000083-55-6	53
4	5-Amino-3-phenylpyrazole	159	C9H9N3	000827-41-8	53
5	benzenamine, 4-(1H-imidazol-1-yl)-	159	C9H9N3	1000404-37-5	50





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Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHB42

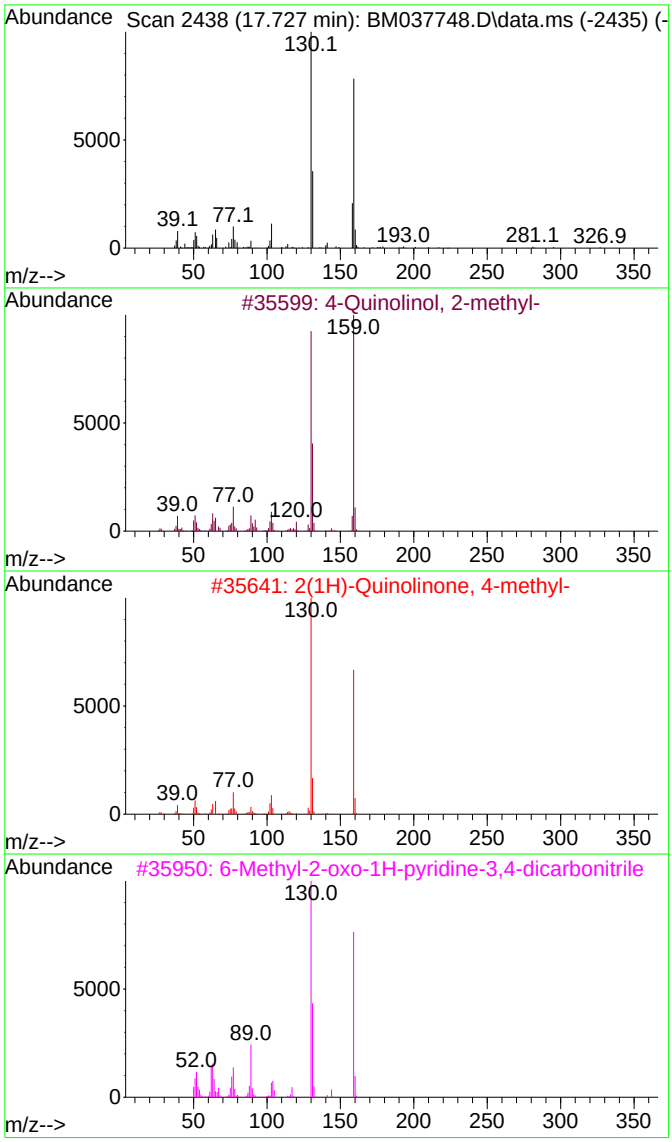
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

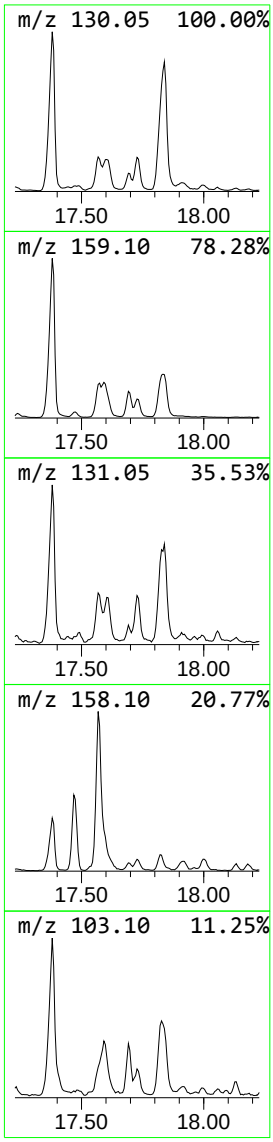
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 4-Quinolinol, 2-methyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.727	2.05 ng/ul	486455	Phenanthrene-d10	17.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Quinolinol, 2-methyl-	159	C10H9NO	000607-67-0	87
2	2(1H)-Quinolinone, 4-methyl-	159	C10H9NO	000607-66-9	72
3	6-Methyl-2-oxo-1H-pyridine-3,4-d...	159	C8H5N3O	100960-56-3	72
4	1-propyl-1H-indole	159	C11H13N	016885-94-2	72
5	6-Methyl-4(1H)-quinolinone	159	C10H9NO	023432-40-8	68





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Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHB42

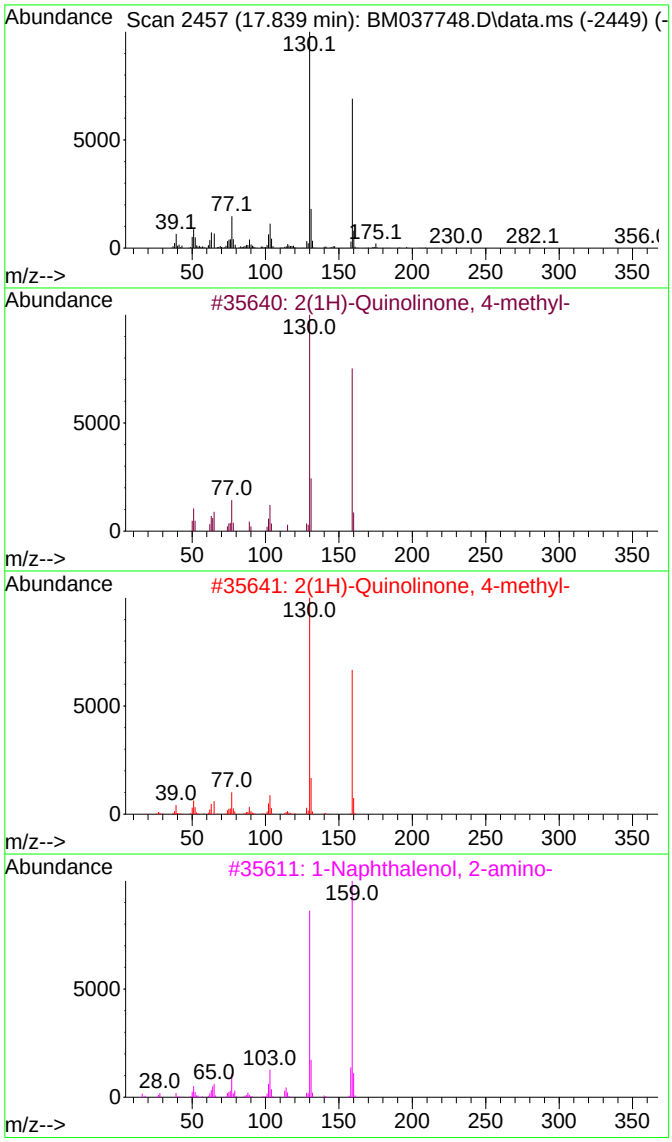
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

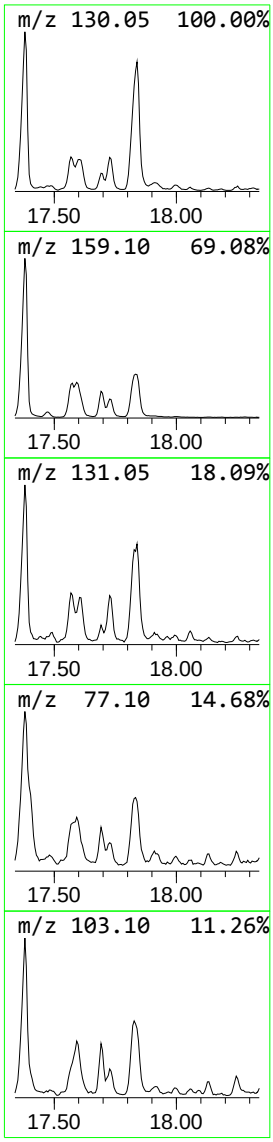
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 2(1H)-Quinolinone, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.839	9.57 ng/ul	2274000	Phenanthrene-d10	17.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Quinolinone, 4-methyl-	159	C10H9NO	000607-66-9	96
2	2(1H)-Quinolinone, 4-methyl-	159	C10H9NO	000607-66-9	95
3	1-Naphthalenol, 2-amino-	159	C10H9NO	000606-41-7	91
4	1-Naphthalenol, 2-amino-	159	C10H9NO	000606-41-7	91
5	2(1H)-Quinolinone, 4-methyl-	159	C10H9NO	000607-66-9	91





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Acq On : 27 Nov 2022 00:13
Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHB42

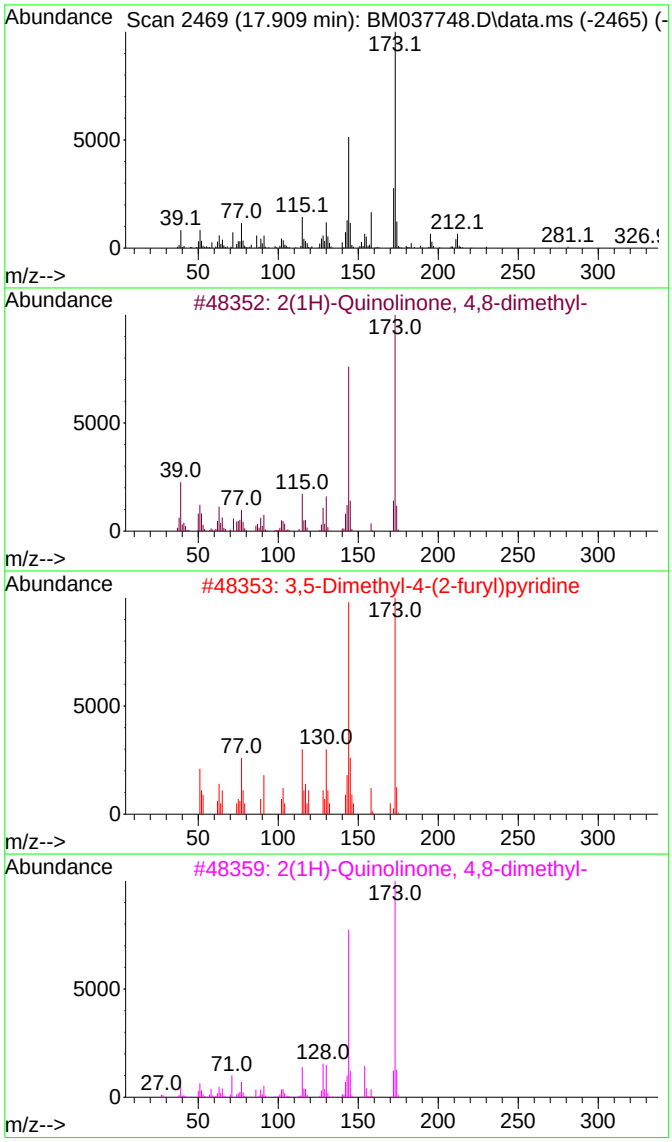
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

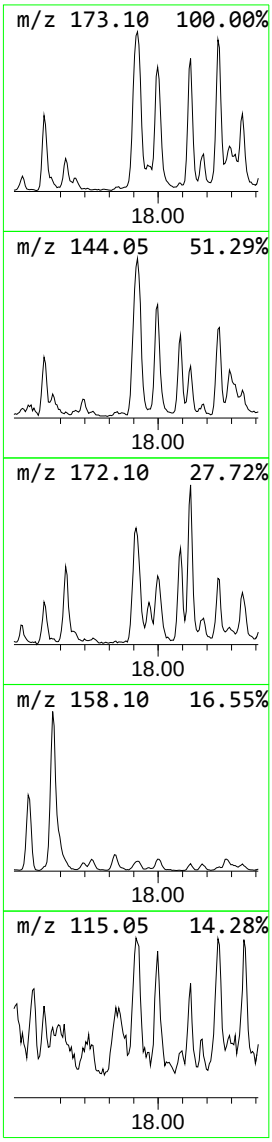
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 2(1H)-Quinolinone, 4,8-dime... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.909	2.25 ng/ul	534476	Phenanthrene-d10	17.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Quinolinone, 4,8-dimethyl-	173	C11H11NO	005349-78-0	93
2	3,5-Dimethyl-4-(2-furyl)pyridine	173	C11H11NO	078563-72-1	90
3	2(1H)-Quinolinone, 4,8-dimethyl-	173	C11H11NO	005349-78-0	87
4	2(1H)-Quinolinone, 4,6-dimethyl-	173	C11H11NO	023947-37-7	83
5	1,6-Dimethyl-2(1H)-quinolinone	173	C11H11NO	029969-49-1	80





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Acq On : 27 Nov 2022 00:13
Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHB42

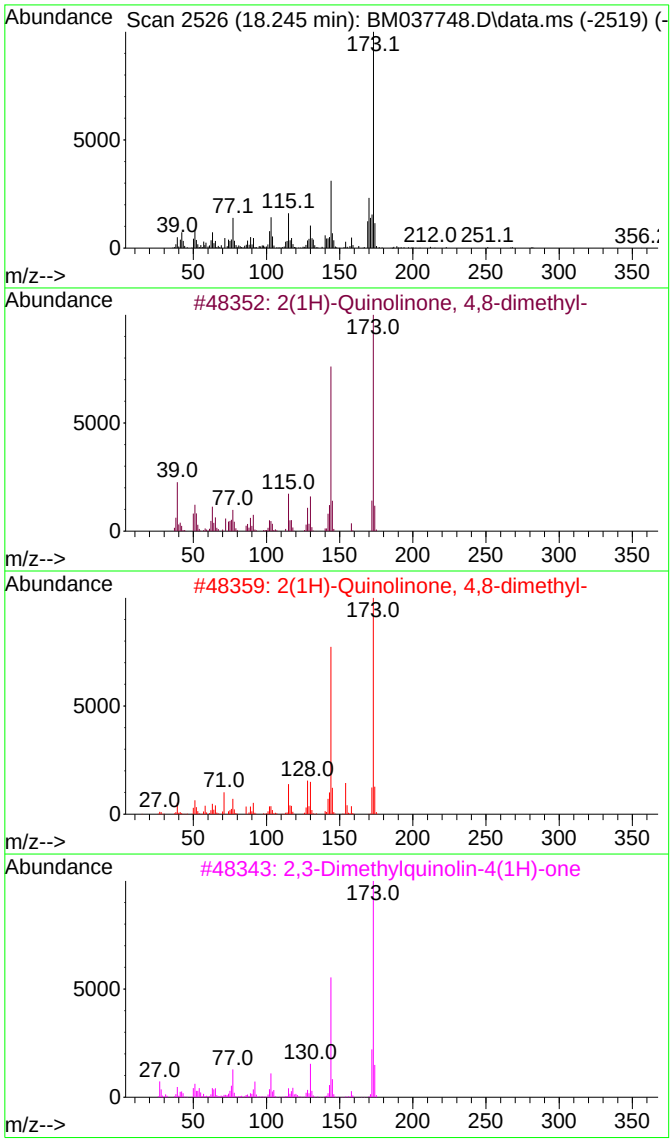
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

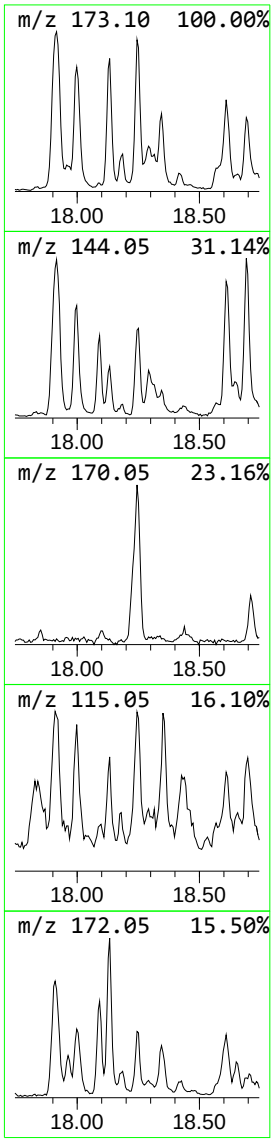
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 2,3-Dimethylquinolin-4(1H)-one Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.245	2.44 ng/ul	581001	Phenanthrene-d10	17.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Quinolinone, 4,8-dimethyl-	173	C11H11NO	005349-78-0	76
2	2(1H)-Quinolinone, 4,8-dimethyl-	173	C11H11NO	005349-78-0	70
3	2,3-Dimethylquinolin-4(1H)-one	173	C11H11NO	058596-45-5	64
4	4-Quinolinol, 2,8-dimethyl-	173	C11H11NO	1010351-37-0	59
5	3-Methyl-4-phenyl-1H-pyrazol-5-a...	173	C10H11N3	060419-81-0	53





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Acq On : 27 Nov 2022 00:13
Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

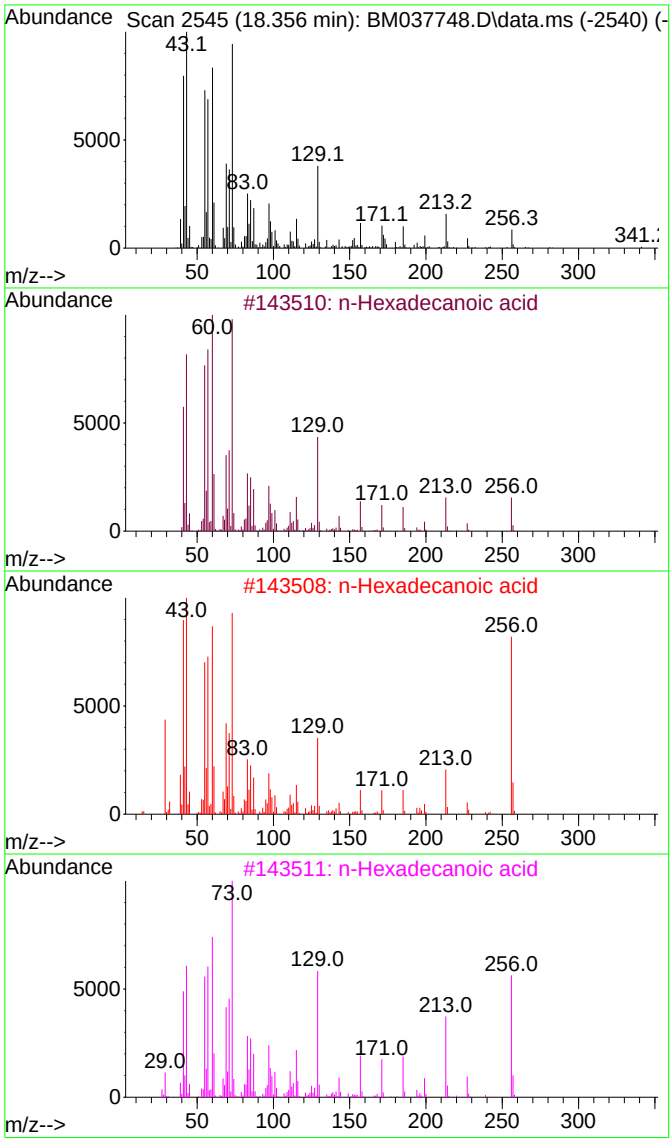
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

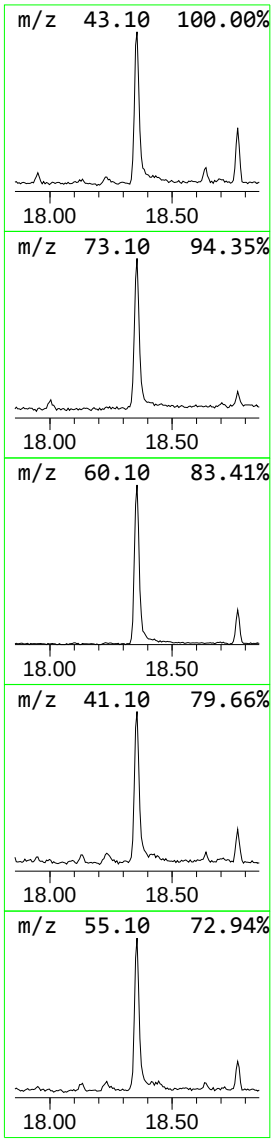
****

Peak Number 23 n-Hexadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.356	3.89 ng/ul	923615	Phenanthrene-d10	17.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	91





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Acq On : 27 Nov 2022 00:13
Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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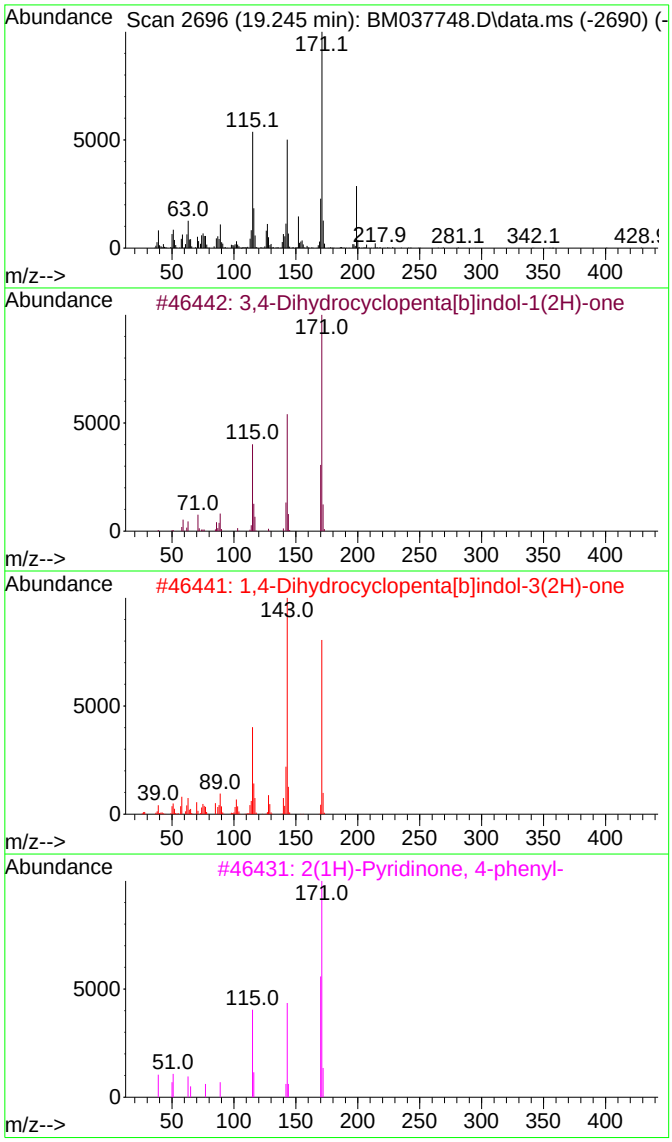
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

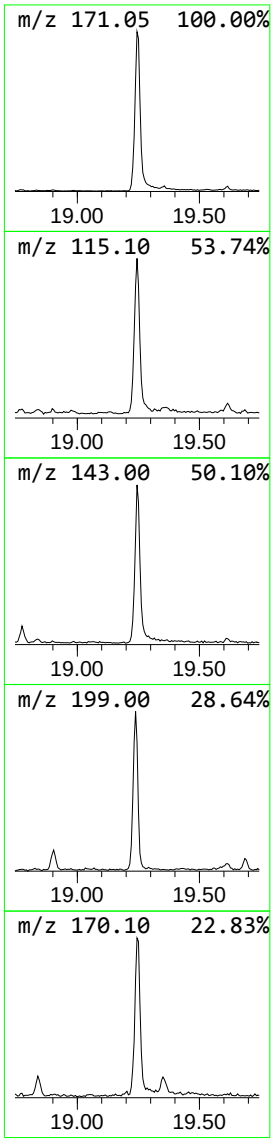
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 3,4-Dihydrocyclopenta[b]ind... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.245	4.55 ng/ul	1082310	Phenanthrene-d10	17.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4-Dihydrocyclopenta[b]indol-1(...	171	C11H9NO	1000305-37-5	76
2	1,4-Dihydrocyclopenta[b]indol-3(...	171	C11H9NO	016244-15-8	70
3	2(1H)-Pyridinone, 4-phenyl-	171	C11H9NO	019006-81-6	64
4	2-(Phenylamino)cyclopenta-2,4-di...	171	C11H9NO	1000436-42-6	60
5	Pyridine, 4-phenyl-, 1-oxide	171	C11H9NO	001131-61-9	60





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Acq On : 27 Nov 2022 00:13
Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHB42

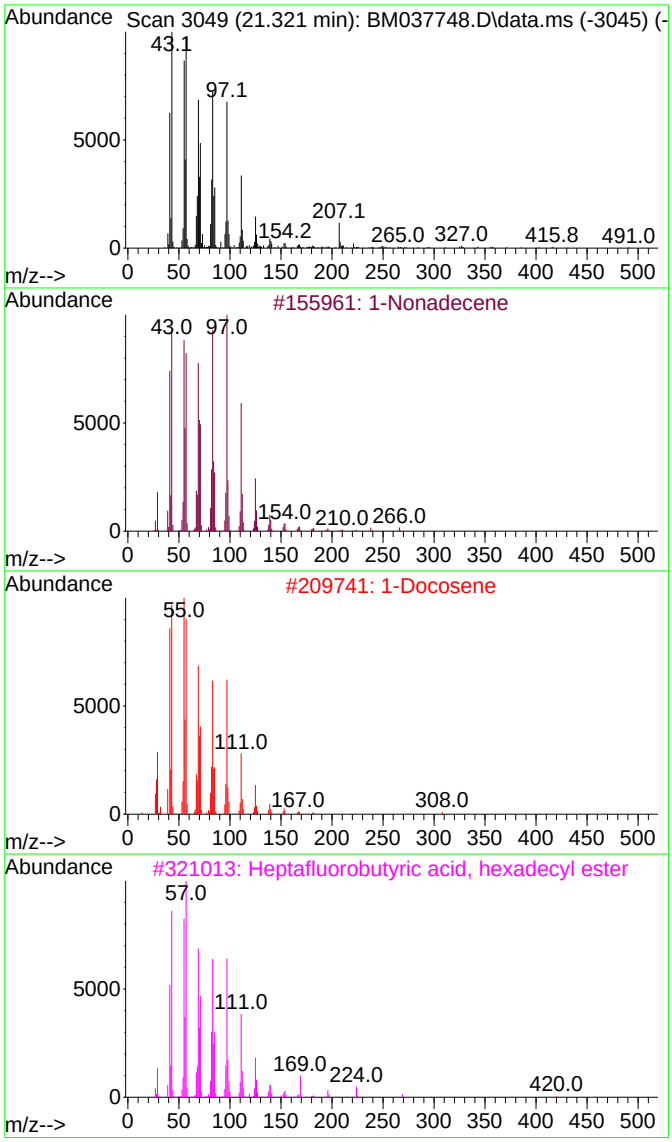
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

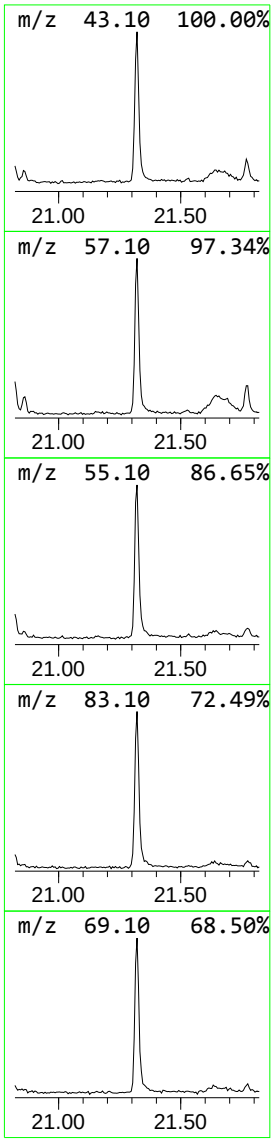
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.321	3.56 ng/ul	786116	Chrysene-d12	21.633

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene	266	C19H38	018435-45-5	96
2	1-Docosene	308	C22H44	001599-67-3	94
3	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91
4	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91
5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91





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Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHB42

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

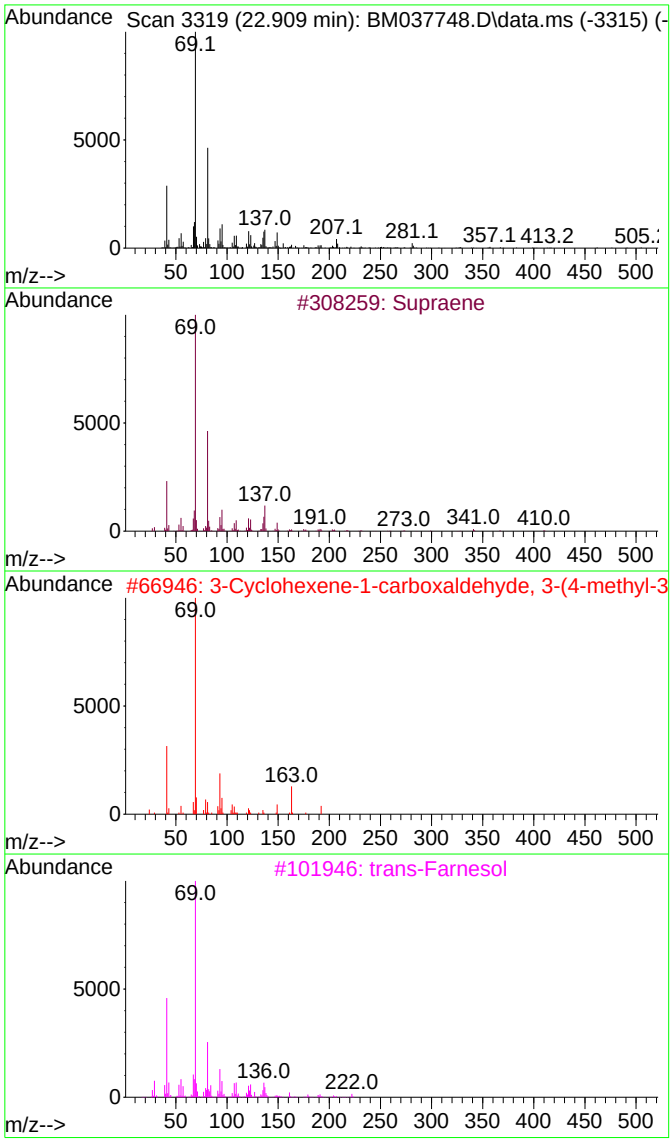
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

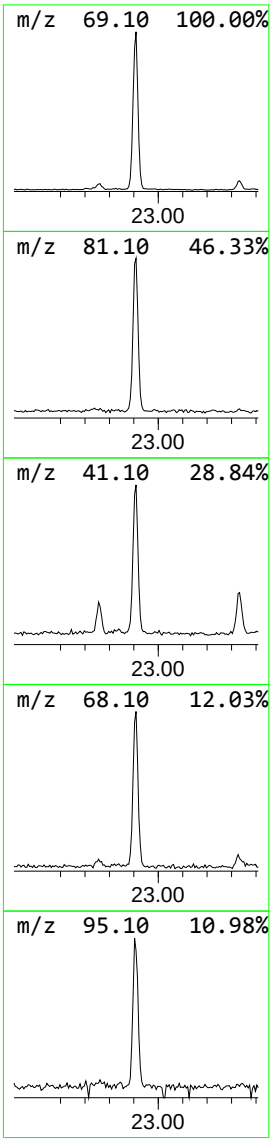
****

Peak Number 26 Supraene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.909	3.19 ng/ul	616160	Perylene-d12	24.127

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Supraene	410	C30H50	007683-64-9	83
2	3-Cyclohexene-1-carboxaldehyde, ...	192	C13H20O	052475-89-5	64
3	trans-Farnesol	222	C15H26O	000106-28-5	64
4	2,6,10,14-Hexadecatetraen-1-ol, ...	332	C22H36O2	061691-98-3	64
5	7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
Data File : BM037748.D
Acq On : 27 Nov 2022 00:13
Operator : CG/JU
Sample : N5694-06
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BHB42

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Indane	8.486	3.1	ng/ul	223643	1	8.110	1462370	20.0
3-Methylphenyla...	8.651	12.5	ng/ul	910801	1	8.110	1462370	20.0
Benzofuran, 2-m...	9.475	36.0	ng/ul	2631070	1	8.110	1462370	20.0
2-Propenal, 3-p...	9.598	7.8	ng/ul	907684	2	10.928	2322030	20.0
3-Phenyl-2-prop...	9.663	9.1	ng/ul	1050880	2	10.928	2322030	20.0
Benzo[b]thiophe...	12.039	10.2	ng/ul	1188590	2	10.928	2322030	20.0
Benzo[b]thiophe...	12.474	4.5	ng/ul	522440	2	10.928	2322030	20.0
Methylchromone	15.127	2.1	ng/ul	329754	3	14.733	3192810	20.0
Indane-5-carbox...	15.210	2.5	ng/ul	406689	3	14.733	3192810	20.0
2-Indolinone, e...	16.080	2.9	ng/ul	459215	3	14.733	3192810	20.0
5-Aminoindan	16.468	39.8	ng/ul	9470120	4	17.468	4754720	20.0
Isoquinoline, 5...	16.651	21.0	ng/ul	4986170	4	17.468	4754720	20.0
benzo[b]thiophe...	16.692	4.7	ng/ul	1117800	4	17.468	4754720	20.0
2(1H)-Quinolino...	17.051	27.5	ng/ul	6528840	4	17.468	4754720	20.0
1-Naphthalenami...	17.233	2.5	ng/ul	598152	4	17.468	4754720	20.0
unknown-01	17.380	66.2	ng/ul	15746400	4	17.468	4754720	20.0
2-Naphthalenol,...	17.692	2.6	ng/ul	610867	4	17.468	4754720	20.0
4-Quinolinol, 2...	17.727	2.0	ng/ul	486455	4	17.468	4754720	20.0
2(1H)-Quinolino...	17.839	9.6	ng/ul	2274000	4	17.468	4754720	20.0
2(1H)-Quinolino...	17.909	2.3	ng/ul	534476	4	17.468	4754720	20.0
2,3-Dimethylqui...	18.245	2.4	ng/ul	581001	4	17.468	4754720	20.0
n-Hexadecanoic ...	18.356	3.9	ng/ul	923615	4	17.468	4754720	20.0
3,4-Dihydrocycl...	19.245	4.5	ng/ul	1082310	4	17.468	4754720	20.0
(DEL) Alkane: S...	21.321	3.6	ng/ul	786116	5	21.633	4421140	20.0
Supraene	22.909	3.2	ng/ul	616160	6	24.127	3859650	20.0