LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BHA80

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Title : SVOA CALIBRATION

Signal : TIC: BP012438.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.969	161	167	184	rVB	11509901	18972643	77.55%	4.765%
2	5.105	348	360	375	rBV	13695529	24464235	100.00%	6.145%
3	5.228	375	381	385	rBV	695721	1002143	4.10%	0.252%
4	5.293	385	392	402	rVB	9177066	14233213	58.18%	3.575%
5	5.422	407	414	423	rVV	4183215	8664044	35.42%	2.176%
6	5.705	451	462	470	rVV	545005	941921	3.85%	0.237%
7	5.793	470	477	488	rVV	3493056	5460553	22.32%	1.371%
8	6.058	513	522	533	rBV2	461746	1147463	4.69%	0.288%
9	6.269	548	558	563	rBV2	1450554	3185242	13.02%	0.800%
10	6.458	583	590	596	rBV2	605827	1072534	4.38%	0.269%
11	6.758	630	641	649	rBV	2194539	3625920	14.82%	0.911%
12	7.046	686	690	696	rVV	642537	975940	3.99%	0.245%
13	7.128	696	704	707	rVV	1022439	1693537	6.92%	0.425%
14	7.169	707	711	716	rVV	2168685	3382989	13.83%	0.850%
15	7.346	732	741	750	rVV2	3598665	7767838	31.75%	1.951%
16	7.428	750	755	761	rVV	2702167	4216734	17.24%	1.059%
17	7.540	769	774	780	rVB	861377	1297825	5.30%	0.326%
18	7.675	789	797	804	rBV4	752057	1705373	6.97%	0.428%
19	7.787	812	816	819	rVV	1010739	1636068	6.69%	0.411%
20	7.822	819	822	829	rVV	1039962	1675512	6.85%	0.421%
21	7.899	829	835	841	rVV	1241107	2152599	8.80%	0.541%
22	7.981	844	849	855	rVB2	777449	1234691	5.05%	0.310%
23	8.157	871	879	884	rVV2	2012408	3810941	15.58%	0.957%
24	8.222	884	890	895	rVV	3370128	5373844	21.97%	1.350%
25	8.275	895	899	904	rVV	1428990	2136720	8.73%	0.537%
26	8.328	904	908	916	rVB	628664	944816	3.86%	0.237%
27	8.416	916	923	934	rBV	4370641	7638629	31.22%	1.919%
28	8.510	934	939	945	rVV	822836	1333293	5.45%	0.335%
29	8.593	945	953	966	rVB3	766466	1957224	8.00%	0.492%
30	8.887	997	1003	1009	rBV2	735160	1247405	5.10%	0.313%
31	8.957	1009	1015	1022	rVV2	1271174	2546449	10.41%	0.640%
32	9.104	1035	1040	1046	rVB	3015532	4665928	19.07%	1.172%
33	9.416	1086	1093	1099	rBV3	582809	1158528	4.74%	0.291%
34	9.675	1131	1137	1150	rVB3	1298722	2985794	12.20%	0.750%
35	9.957	1180	1185	1188	rBV	835774	1481137	6.05%	0.372%
36	10.216	1221	1229	1236	rBV	1650120	3006187	12.29%	0.755%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Title : SVOA CALIBRATION

37	10.593	1286	1293	1297	rBV	1348135	2311947	9.45%	0.581%
38	10.734	1311	1317	1326	rVB3	1376582	3418860	13.97%	0.859%
39	11.004	1359	1363	1369	rVB	921203	1455740	5.95%	0.366%
40	11.187	1388	1394	1403	rVV2	553203	1093938	4.47%	0.275%
41	11.293	1403	1412	1417	rVV2	784665	1613484	6.60%	0.405%
42	11.946	1513	1523	1529	rBV	2473256	4141410	16.93%	1.040%
43	12.993	1691	1701	1706	rVV2	4314236	7561727	30.91%	1.899%
44	13.857	1842	1848	1851	rVV	3093682	4566056	18.66%	1.147%
45	13.887	1851	1853	1858	rVV	1362127	1811581	7.41%	0.455%
46	13.981	1858	1869	1872	rVV	4523158	11906019	48.67%	2.990%
47	14.010	1872	1874	1879	rVV	2812250	3408611	13.93%	0.856%
48	14.092	1883	1888	1891	rVV	1847705	2621139	10.71%	0.658%
49	14.134	1891	1895	1901	rVB	3509619	4973695	20.33%	1.249%
50	14.234	1906	1912	1917	rVV	1051083	1720615	7.03%	0.432%
51	14.469	1941	1952	1962	rVB2	12643036	23268617	95.11%	5.844%
52	14.816	2005	2011	2015	rBV	1222171	1997427	8.16%	0.502%
53	15.351	2097	2102	2106	rBV	921082	1246348	5.09%	0.313%
54	15.445	2112	2118	2125	rVB2	7901613	14066092	57.50%	3.533%
55	15.698	2153	2161	2166	rBV2	1323554	2574908	10.53%	0.647%
56	15.963	2201	2206	2221	rVB3	1401983	2909359	11.89%	0.731%
57	16.086	2221	2227	2232	rBV	1205236	2029296	8.29%	0.510%
58	16.163	2233	2240	2246	rVV2	1430942	2918487	11.93%	0.733%
59	16.381	2272	2277	2286	rVB	3185748	5180973	21.18%	1.301%
60	16.504	2291	2298	2302	rBV6	951835	1917942	7.84%	0.482%
61	16.545	2302	2305	2310	rVB	829355	1056126	4.32%	0.265%
62	16.604	2310	2315	2323	rVB6	622524	1209542	4.94%	0.304%
63	16.804	2344	2349	2354	rBV	4336454	5936670	24.27%	1.491%
64	16.951	2368	2374	2380	rBV	2879718	4987266	20.39%	1.253%
65	17.204	2411	2417	2423	rVB	2945800	4432281	18.12%	1.113%
66	17.304	2428	2434	2439	rBV2	5299263	7920093	32.37%	1.989%
67	17.363	2439	2444	2448	rBV	2541261	3439359	14.06%	0.864%
68	17.698	2497	2501	2512	rBV	735774	1277725	5.22%	0.321%
69	17.904	2532	2536	2539	rVB	873387	1125852	4.60%	0.283%
70	18.098	2565	2569	2575	rBV	1810828	2327539	9.51%	0.585%
71	18.245	2589	2594	2599	rVB2	2924196	4070592	16.64%	1.022%
72	18.322	2603	2607	2609	rBV	1471266	2057673	8.41%	0.517%
73	18.345	2609	2611	2614	rVV	2371673	2978653	12.18%	0.748%
74	18.398	2614	2620	2624	rVB	2428554	4738204	19.37%	1.190%
75	18.639	2657	2661	2666	rBV2	645979	947105	3.87%	0.238%
76	18.763	2677	2682	2688	rBV3	1243069	2205535	9.02%	0.554%

LSC Area Percent Report

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Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Title : SVOA CALIBRATION

77	18.886	2700	2703	2706	rBV	630403	908993	3.72%	0.228%
78	19.122	2739	2743	2750	rBV	2010172	2531701	10.35%	0.636%
79	19.269	2765	2768	2775	rVB2	1132230	1387996	5.67%	0.349%
80	19.333	2775	2779	2789	rVB	1517833	2258237	9.23%	0.567%
81	19.545	2810	2815	2820	rBV	6038387	8075969	33.01%	2.028%
82	19.598	2820	2824	2834	rVV	2456558	3888656	15.90%	0.977%
83	19.845	2862	2866	2873	rVB	919124	1440517	5.89%	0.362%
84	19.998	2888	2892	2895	rBV	991206	1160554	4.74%	0.291%
85	20.045	2895	2900	2909	rBV3	2706388	6868257	28.07%	1.725%
86	20.127	2910	2914	2919	rVB2	1514629	2133703	8.72%	0.536%
87	20.398	2957	2960	2965	rVB2	722012	937694	3.83%	0.236%
88	20.539	2981	2984	2992	rVB3	1093758	1586107	6.48%	0.398%
89	20.998	3057	3062	3071	rVB	6740626	8951539	36.59%	2.248%
90	21.298	3109	3113	3123	rVB	3757224	5249946	21.46%	1.319%
91	21.421	3131	3134	3138	rBV	1473171	1976049	8.08%	0.496%
92	21.469	3138	3142	3146	rVB2	4354558	5552892	22.70%	1.395%
93	21.521	3146	3151	3155	rBV	5618091	7197405	29.42%	1.808%
94	21.569	3155	3159	3161	rVV	2067028	2615278	10.69%	0.657%
95	21.598	3161	3164	3169	rVB	2255608	2716743	11.10%	0.682%
96	21.957	3222	3225	3230	rVB2	856893	1089300	4.45%	0.274%
97	22.051	3238	3241	3250	rVB	789734	1136503	4.65%	0.285%
98	23.533	3486	3493	3506	rVB2	4104222	8691058	35.53%	2.183%
99	23.686	3513	3519	3528	rVB2	2191552	4425406	18.09%	1.112%
100	25.080	3750	3756	3770	rVB6	1012761	3144528	12.85%	0.790%

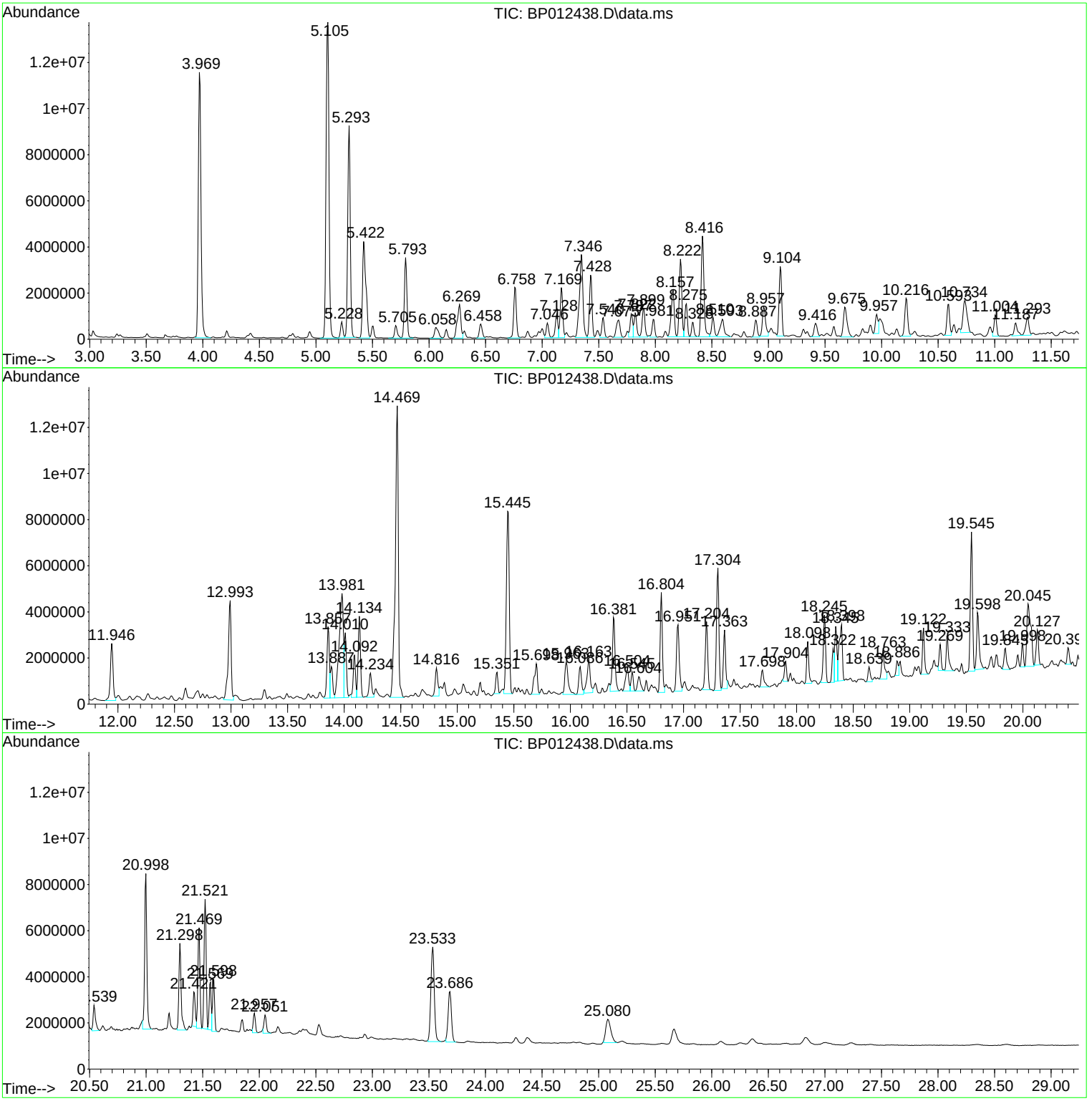
Sum of corrected areas: 398145429

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

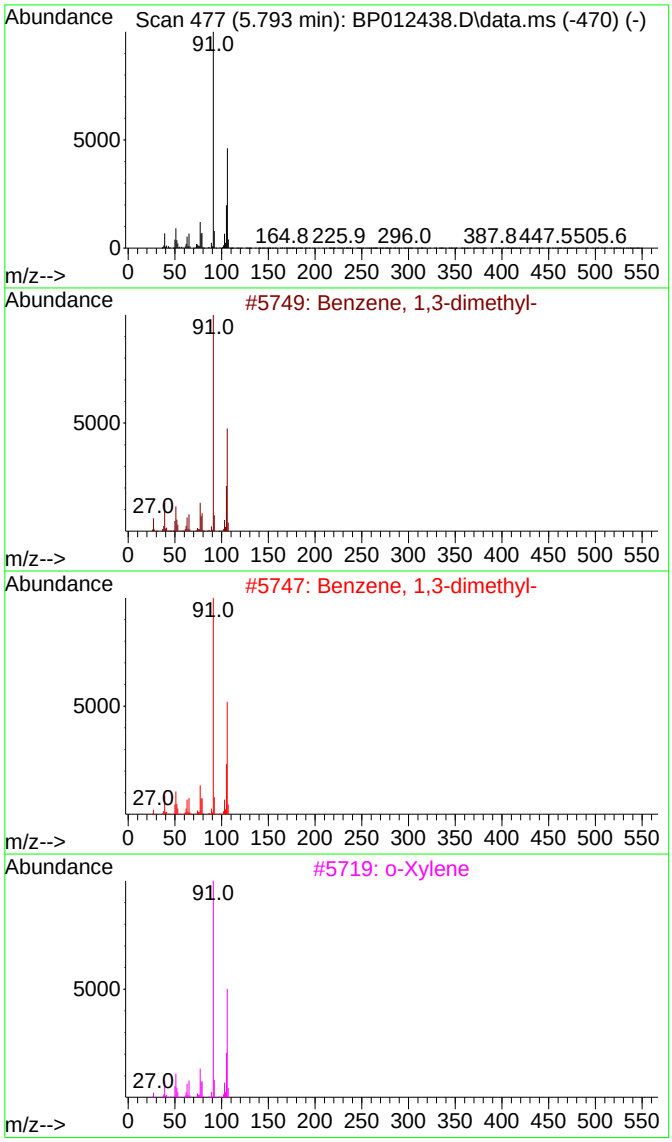
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

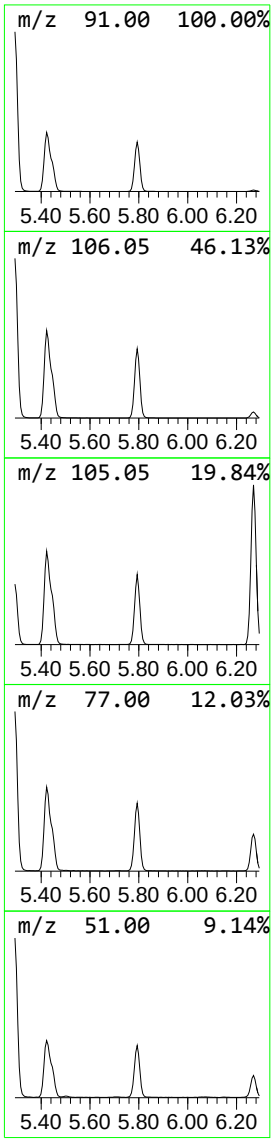
****

Peak Number 7 Benzene, 1,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.793	66.75 ng/ul	5460550	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
3	o-Xylene	106	C8H10	000095-47-6	97
4	p-Xylene	106	C8H10	000106-42-3	97
5	p-Xylene	106	C8H10	000106-42-3	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

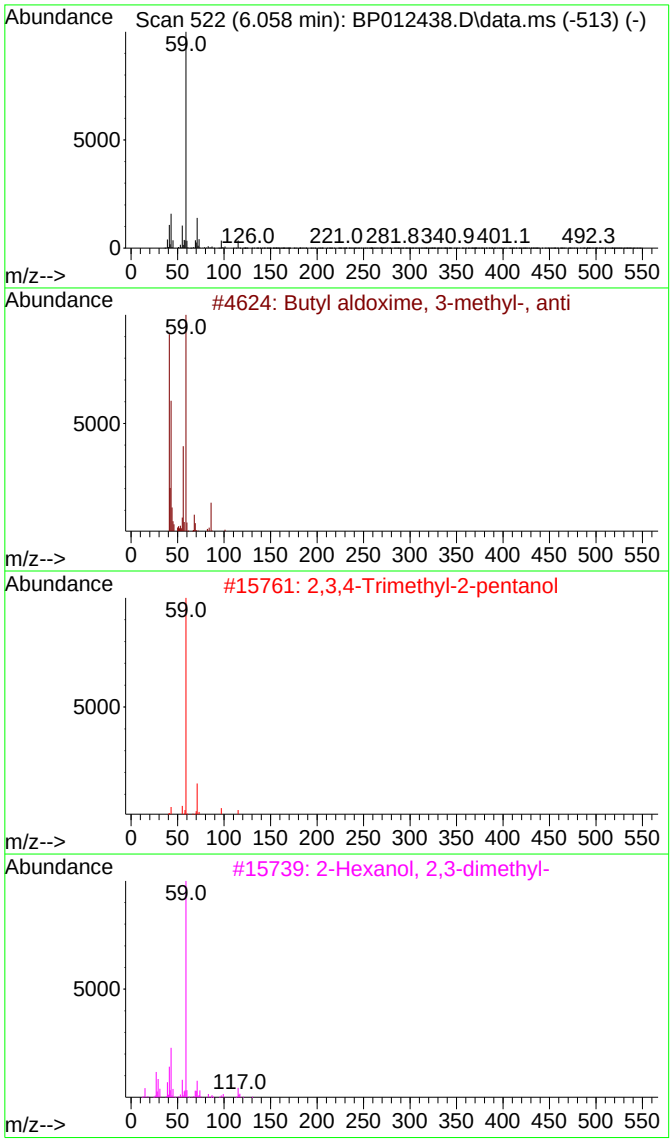
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

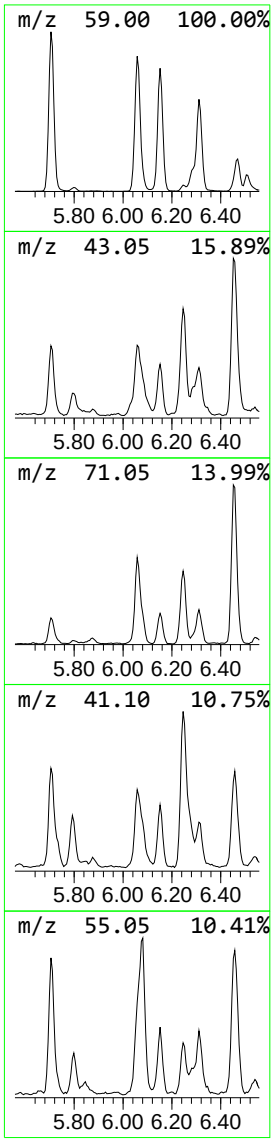
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Butyl aldoxime, 3-methyl-, ... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.058	14.03 ng/ul	1147460	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl aldoxime, 3-methyl-, anti	101	C5H11NO	005775-74-6	59
2	2,3,4-Trimethyl-2-pentanol	130	C8H18O	066576-26-9	50
3	2-Hexanol, 2,3-dimethyl-	130	C8H18O	019550-03-9	50
4	6-Methyl-2-heptanol, methyl ether	144	C9H20O	1000365-52-0	50
5	2-Decanol, methyl ether	172	C11H24O	1000333-83-9	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

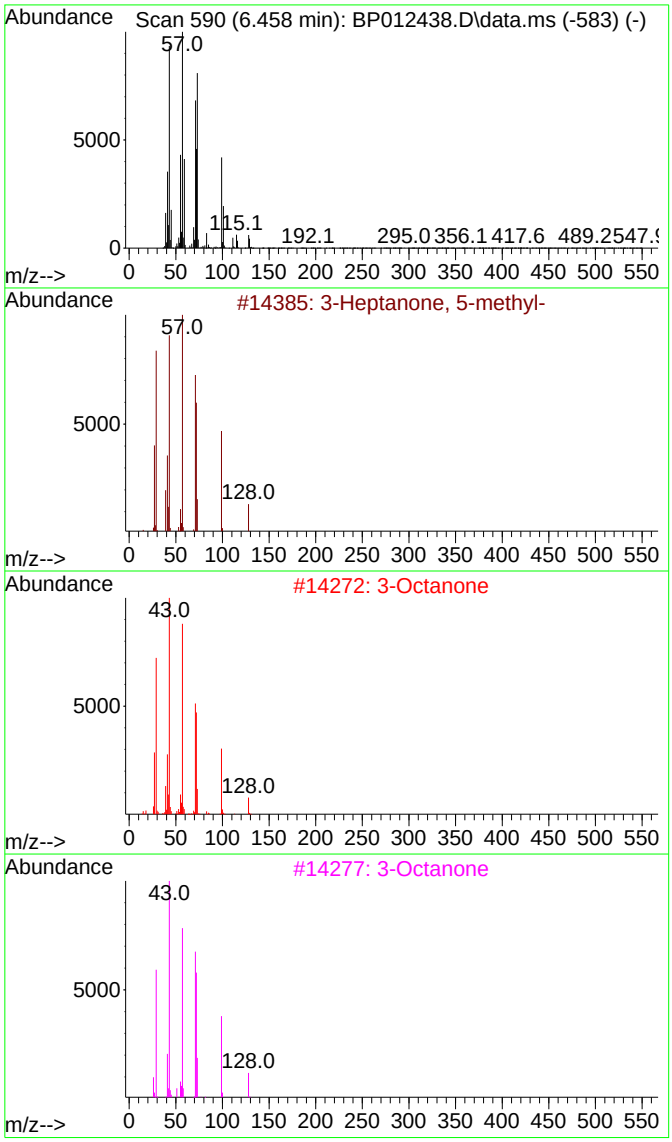
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

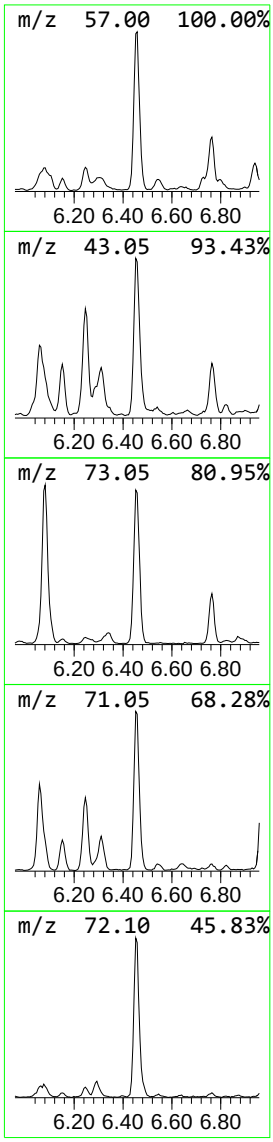
****

Peak Number 10 unknown-01 Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.458	13.11 ng/ul	1072530	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	47
2	3-Octanone	128	C8H16O	000106-68-3	46
3	3-Octanone	128	C8H16O	000106-68-3	43
4	Ethyl amylketone	128	C8H16O	020616-93-7	43
5	Hexaborane	76	B6H10	023777-80-2	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
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Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

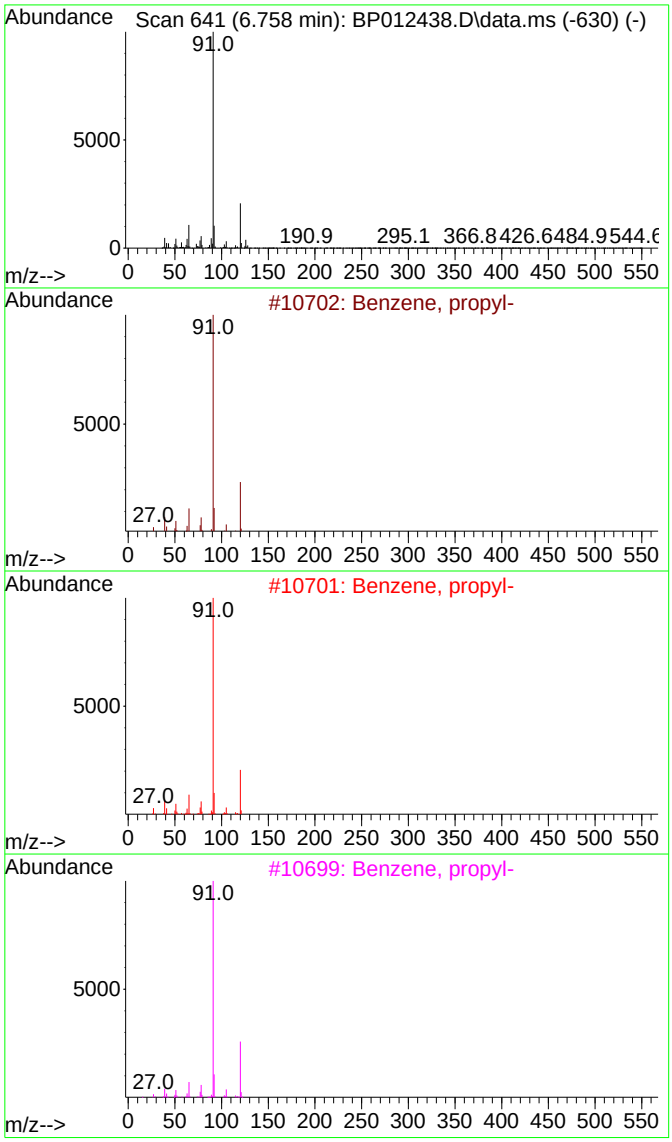
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

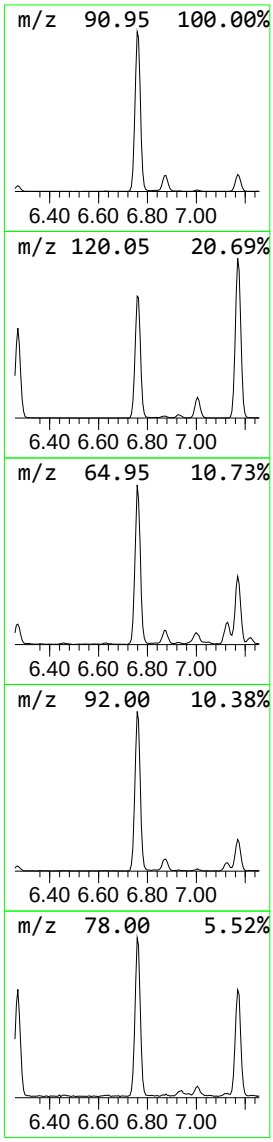
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, propyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.758	44.32 ng/ul	3625920	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	87
2	Benzene, propyl-	120	C9H12	000103-65-1	81
3	Benzene, propyl-	120	C9H12	000103-65-1	74
4	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72
5	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

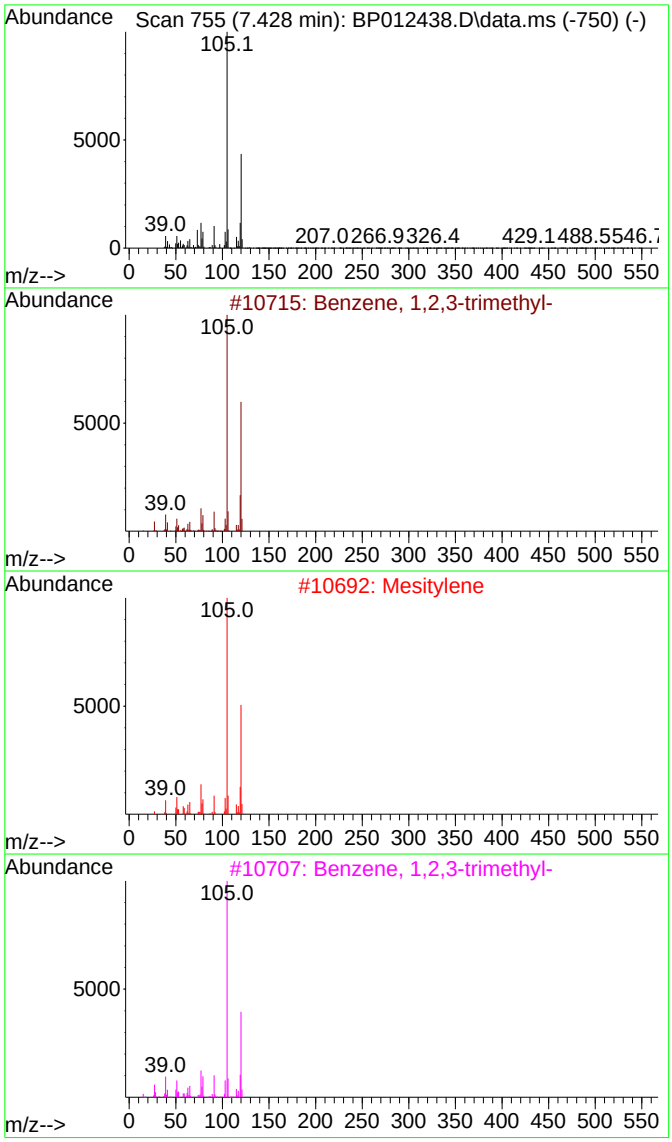
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

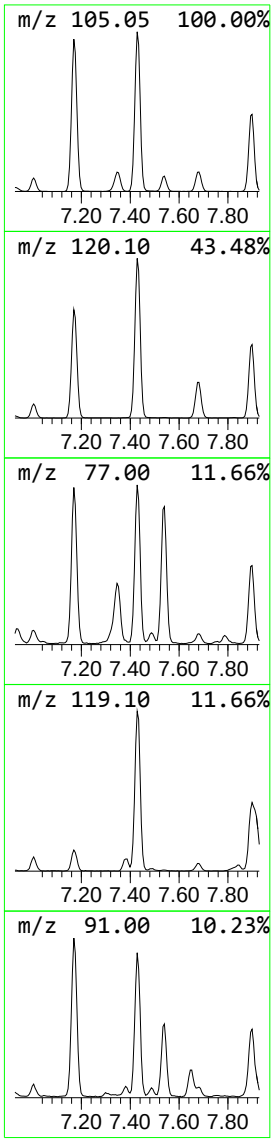
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, 1,2,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.428	51.55 ng/ul	4216730	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Mesitylene	120	C9H12	000108-67-8	97
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4	Mesitylene	120	C9H12	000108-67-8	94
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94





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Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

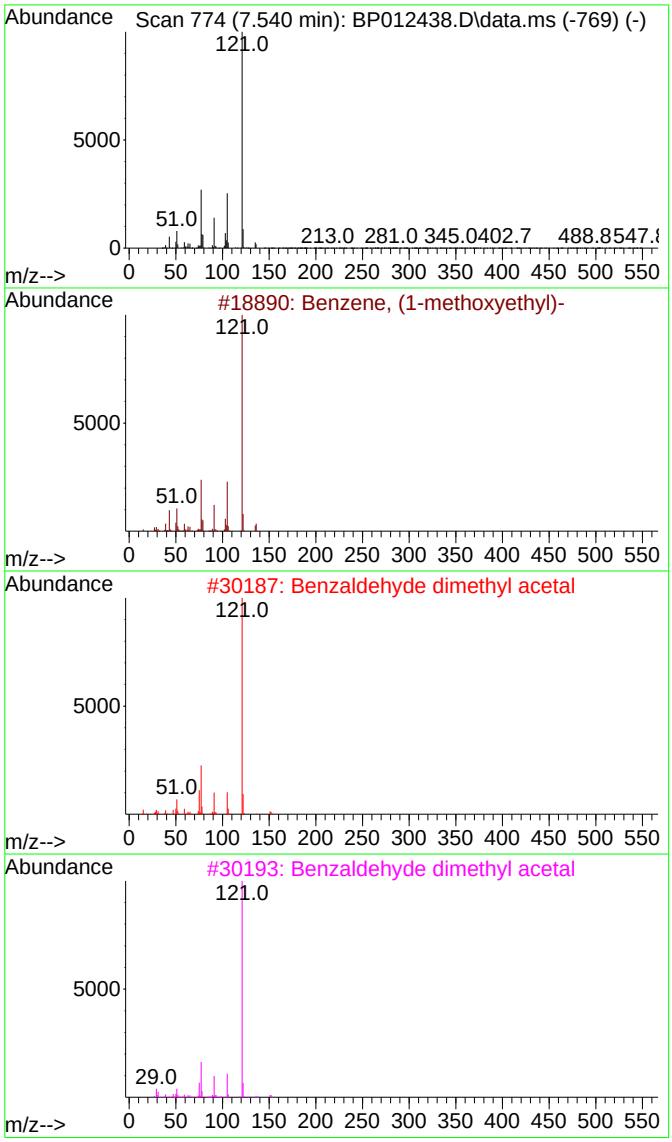
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

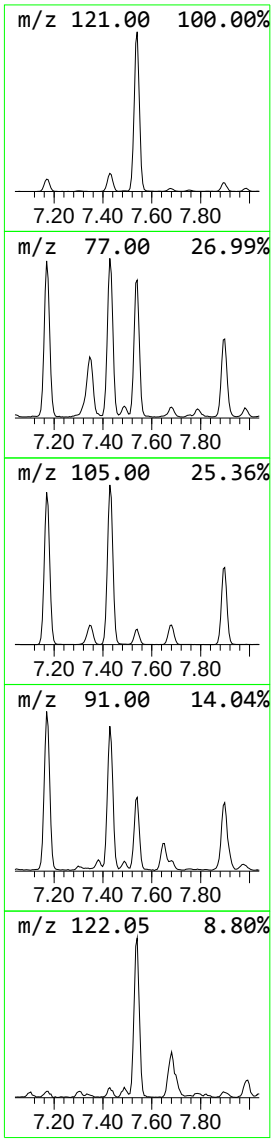
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, (1-methoxyethyl)- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.540	15.87 ng/ul	1297830	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methoxyethyl)-	136	C9H12O	004013-34-7	94
2	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	64
3	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	64
4	Benzenamine, 3,4-dimethyl-	121	C8H11N	000095-64-7	64
5	Benzene, 1-ethyl-4-methoxy-	136	C9H12O	001515-95-3	59





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Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

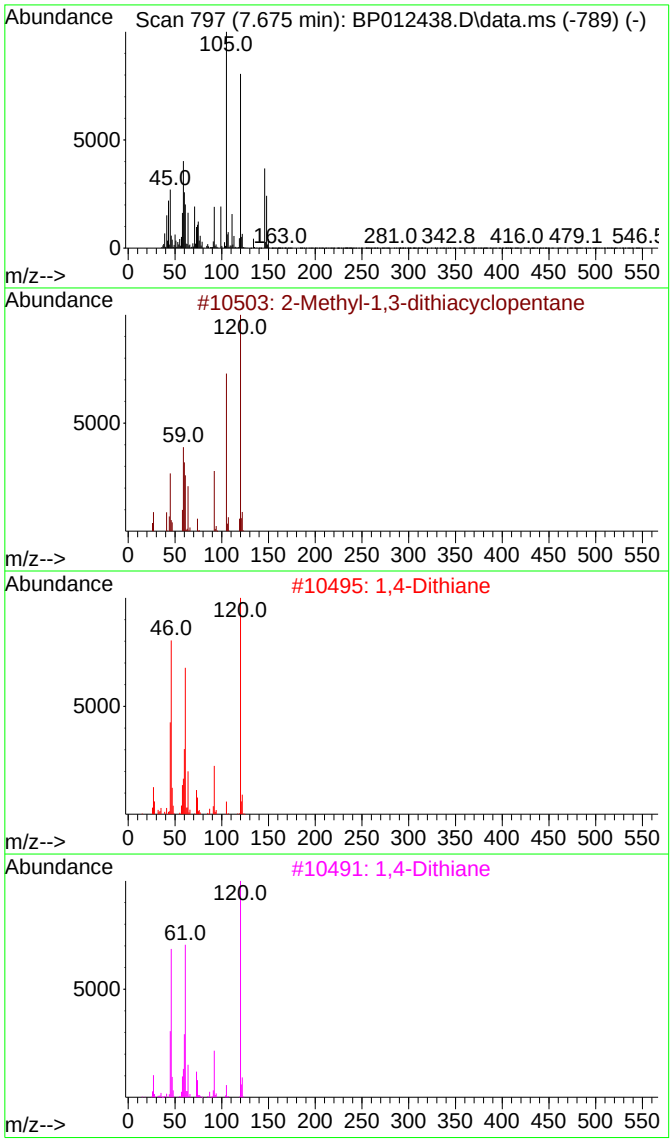
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

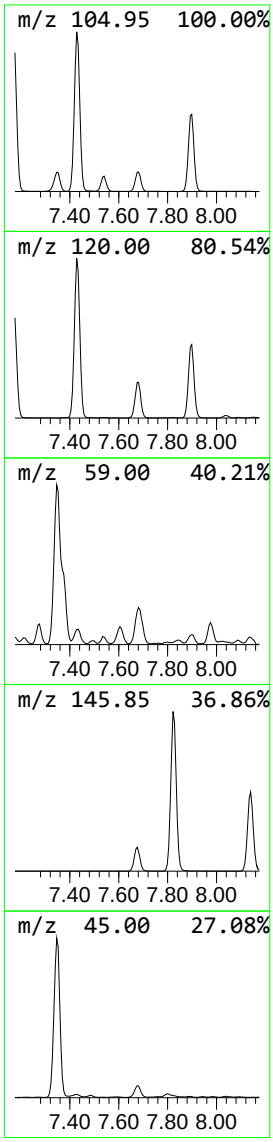
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 2-Methyl-1,3-dithiacyclopen... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.675	20.85 ng/ul	1705370	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-1,3-dithiacyclopentane	120	C4H8S2	005616-51-3	91
2	1,4-Dithiane	120	C4H8S2	000505-29-3	38
3	1,4-Dithiane	120	C4H8S2	000505-29-3	35
4	Ethanedithioamide	120	C2H4N2S2	000079-40-3	35
5	1,4-Dithiane	120	C4H8S2	000505-29-3	25





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Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

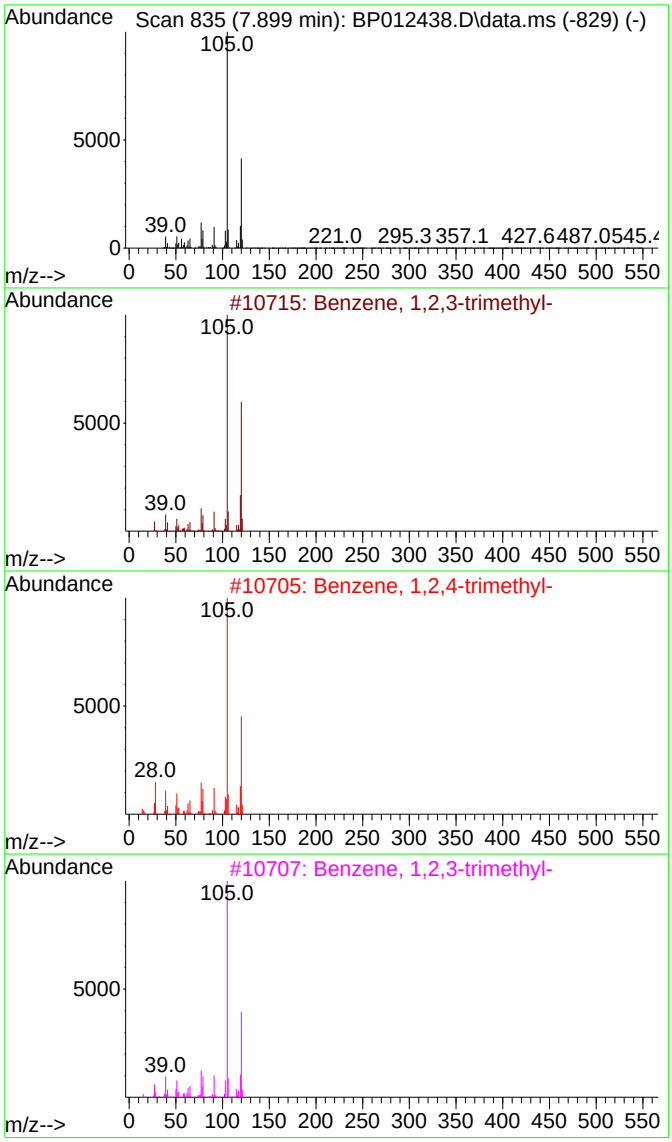
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

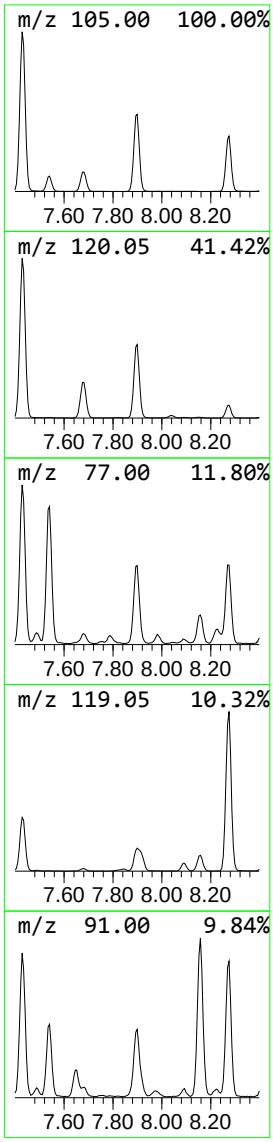
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzene, 1,2,4-trimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.899	26.31 ng/ul	2152600	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Mesitylene	120	C9H12	000108-67-8	94
5	Mesitylene	120	C9H12	000108-67-8	93





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Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

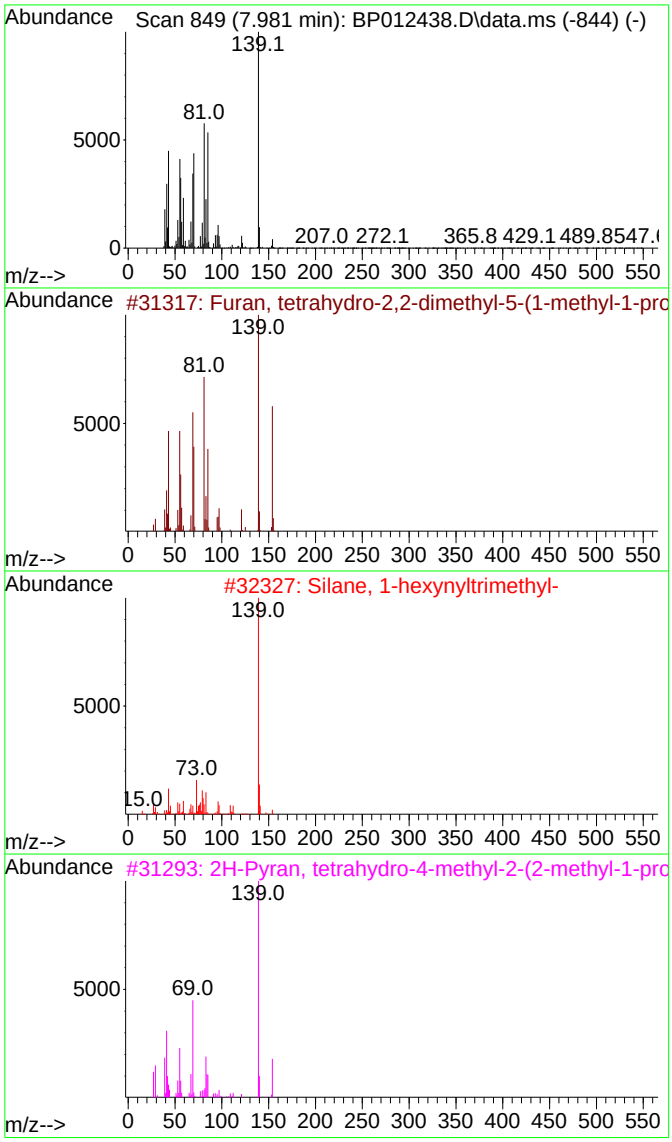
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

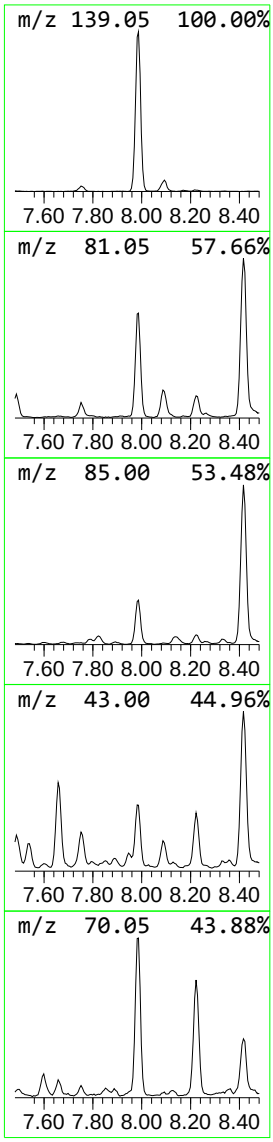
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Furan, tetrahydro-2,2-dimet... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.981	15.09 ng/ul	1234690	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Furan, tetrahydro-2,2-dimethyl-5...	154	C10H18O	007416-35-5	56
2	Silane, 1-hexynyltrimethyl-	154	C9H18Si	003844-94-8	27
3	2H-Pyran, tetrahydro-4-methyl-2-...	154	C10H18O	016409-43-1	27
4	2-Fluoro-.alpha.-methylbenzyl al...	154	C9H11FO	1000365-11-2	27
5	trans-Rose oxide	154	C10H18O	000876-18-6	25





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Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

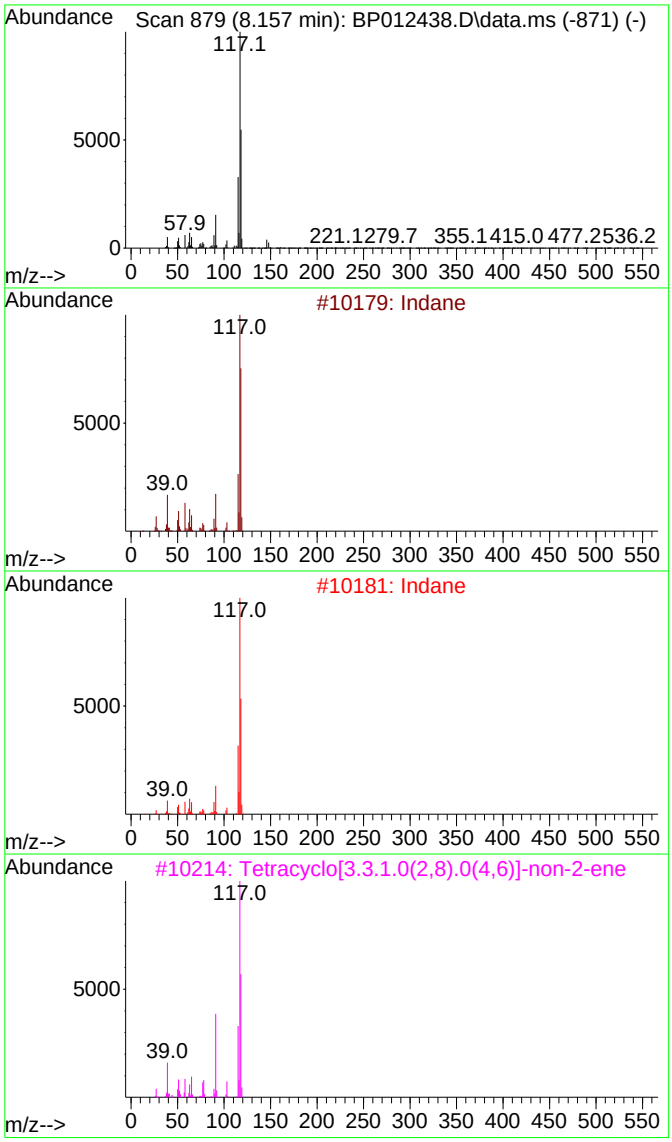
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

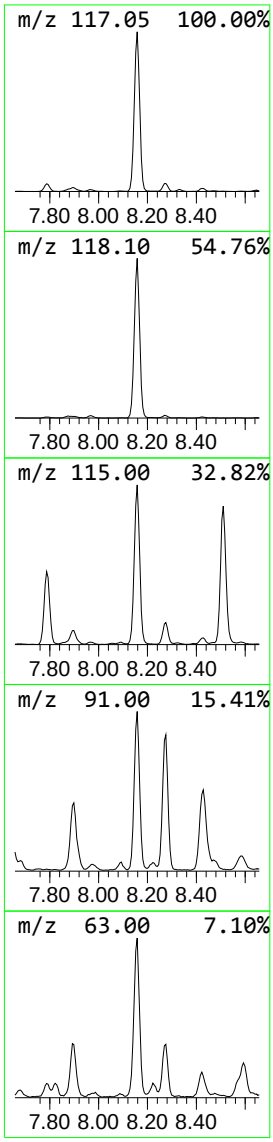
****

Peak Number 17 Indane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.157	46.59 ng/ul	3810940	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Indane	118	C9H10	000496-11-7	81
3	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
4	Indane	118	C9H10	000496-11-7	60
5	Benzene, 1-(chloromethyl)-4-ethe...	152	C9H9Cl	001592-20-7	59





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Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

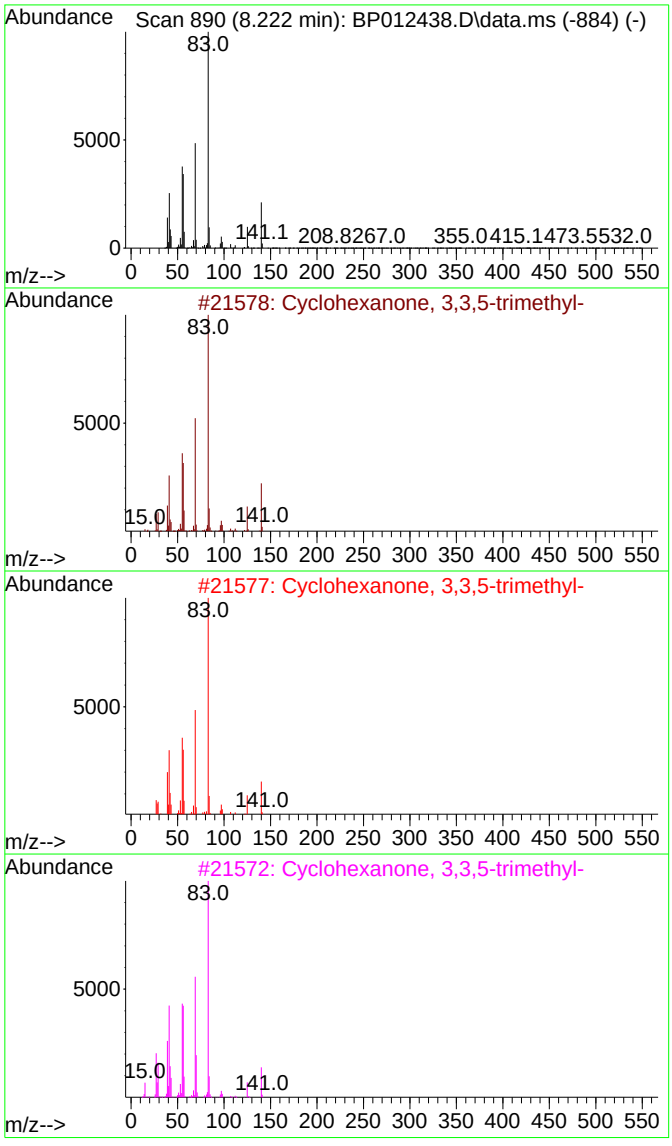
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

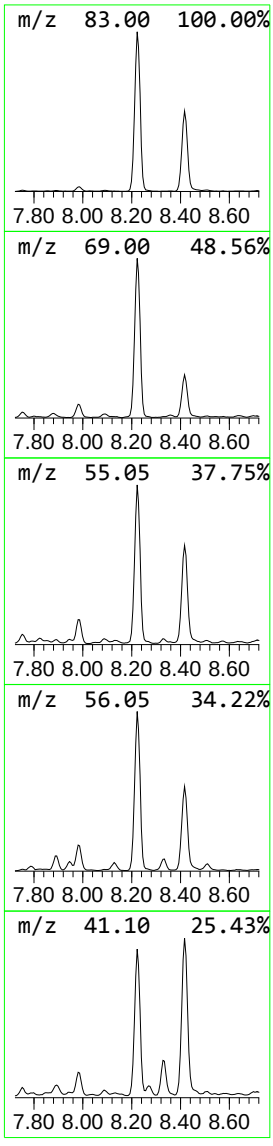
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Cyclohexanone, 3,3,5-trimet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.222	65.69 ng/ul	5373840	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	95
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	94
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
4	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
5	Cyclohexane, butyl-	140	C10H20	001678-93-9	59





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Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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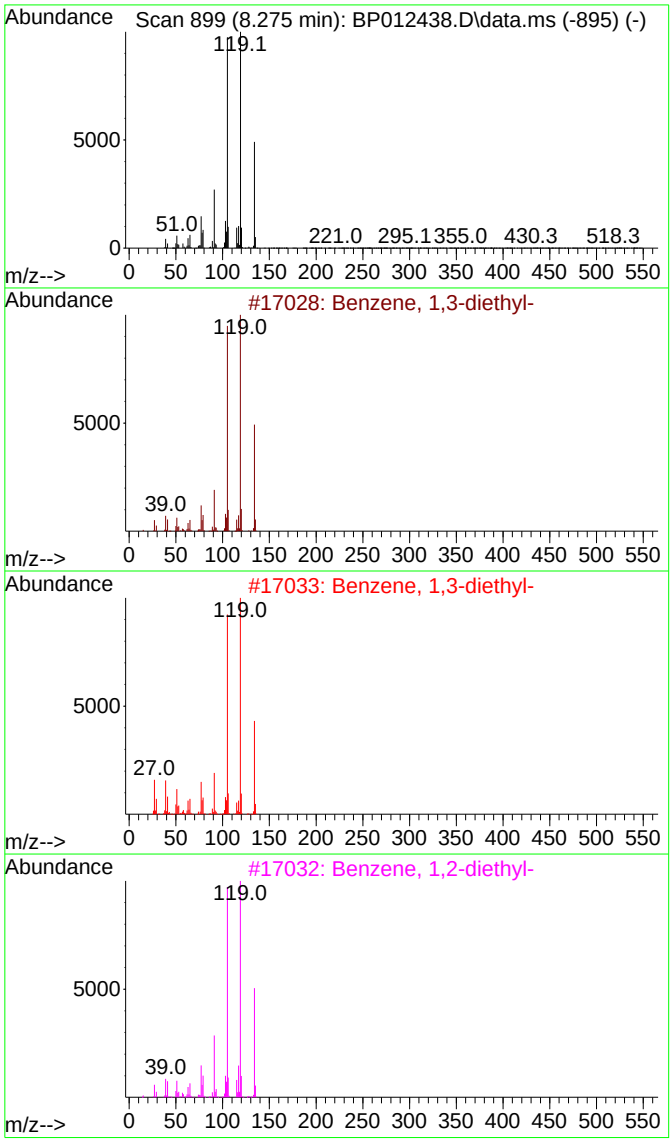
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

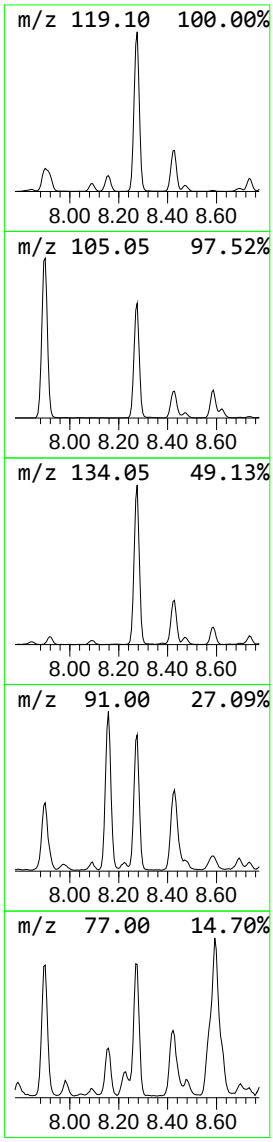
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Benzene, 1,3-diethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.275	26.12 ng/ul	2136720	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
2	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
3	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	94
4	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94





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Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

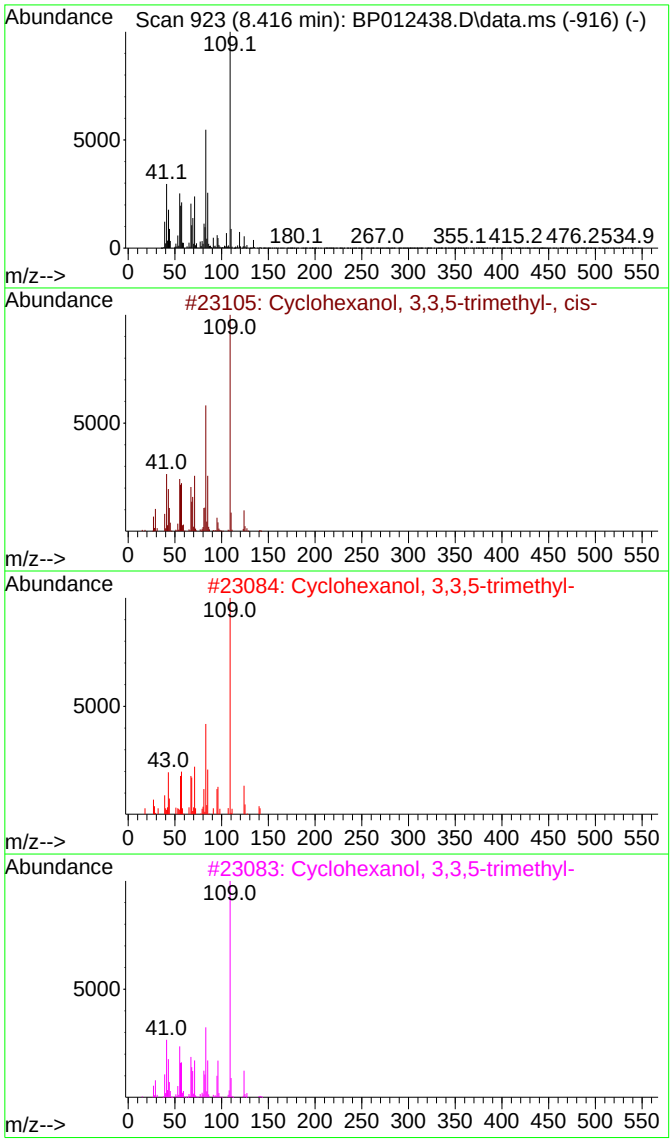
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

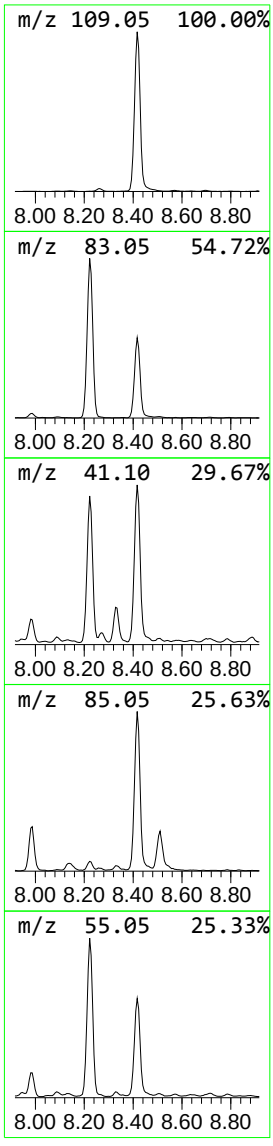
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Cyclohexanol, 3,3,5-trimeth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.416	93.38 ng/ul	7638630	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanol, 3,3,5-trimethyl-, ...	142	C9H18O	000933-48-2	86
2	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	78
3	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	45
4	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	35
5	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	35





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Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

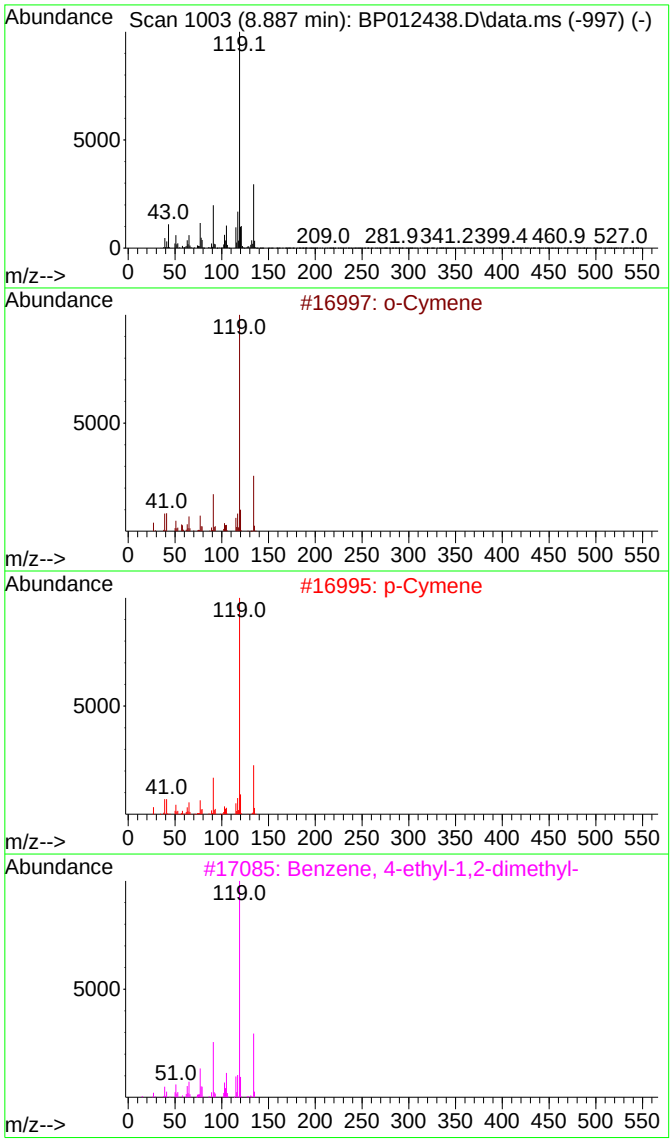
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

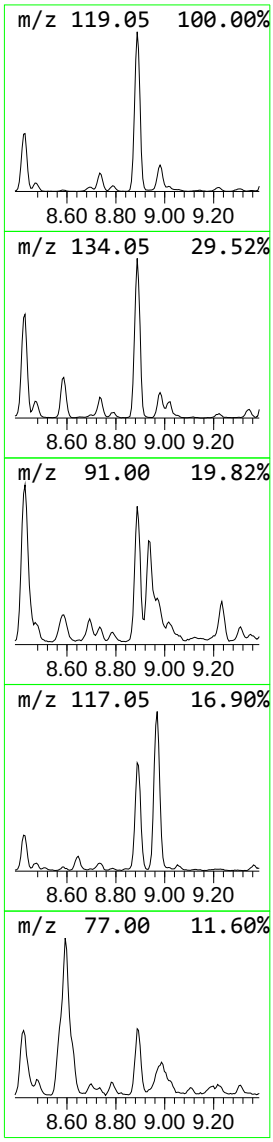
****

Peak Number 22 o-Cymene Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.887	15.25 ng/ul	1247410	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	95
2	p-Cymene	134	C10H14	000099-87-6	94
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
4	o-Cymene	134	C10H14	000527-84-4	93
5	o-Cymene	134	C10H14	000527-84-4	93





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Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

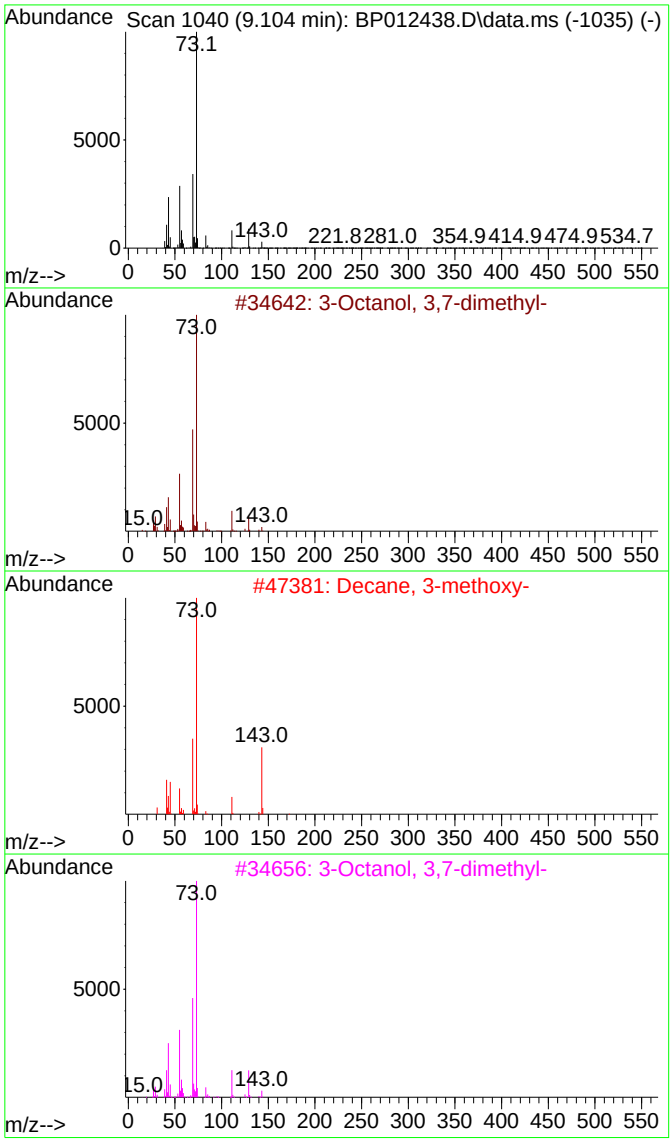
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

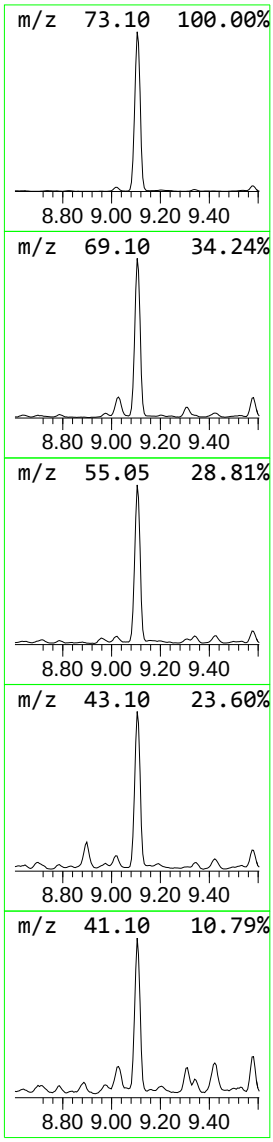
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 3-Octanol, 3,7-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.104	57.04 ng/ul	4665930	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	64
2	Decane, 3-methoxy-	172	C11H24O	055955-64-1	59
3	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	53
4	3-Octanol, 3,6-dimethyl-	158	C10H22O	000151-19-9	42
5	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	42





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Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

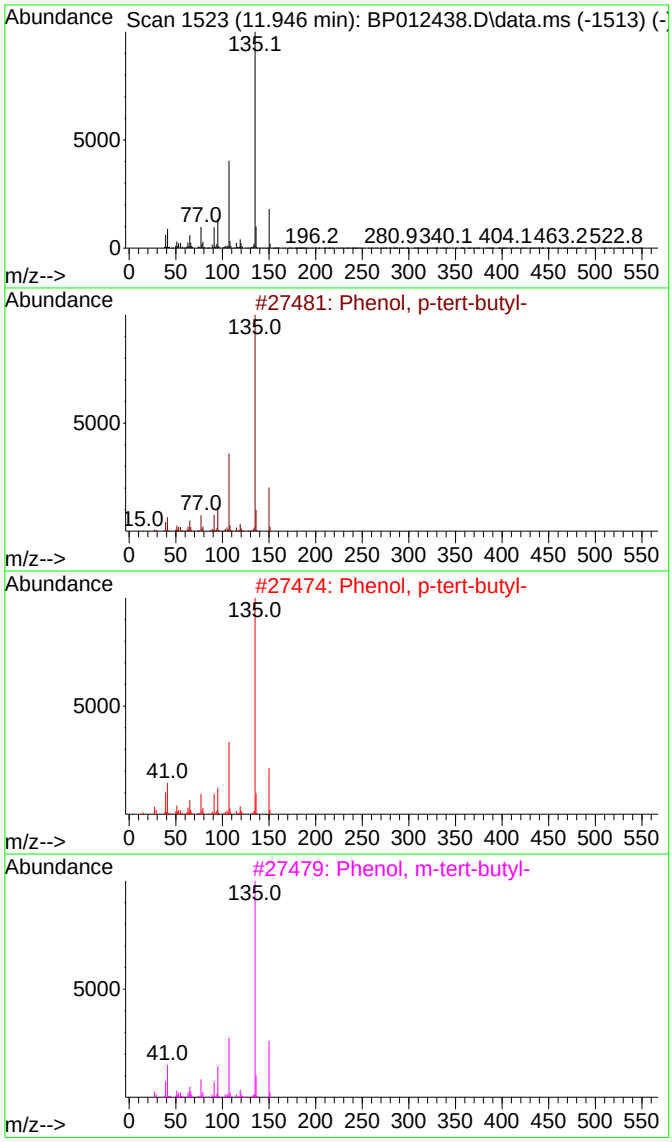
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

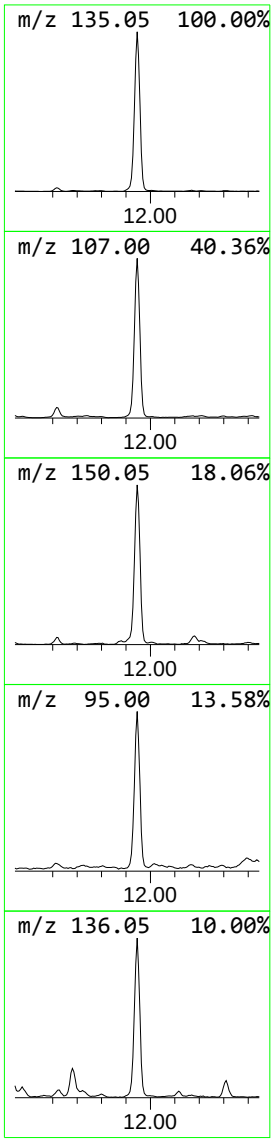
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 Phenol, p-tert-butyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.945	35.83 ng/ul	4141410	Naphthalene-d8	10.593

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	97
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	96
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	94
4	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	87





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Operator : CG/JU
Sample : N5300-03
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ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

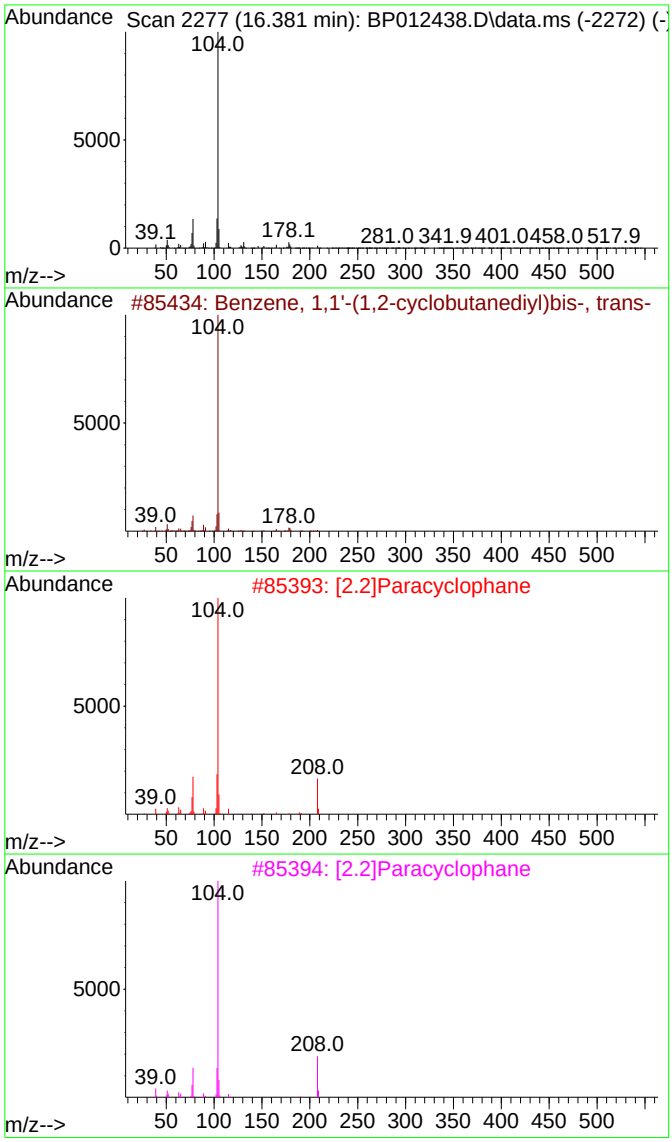
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

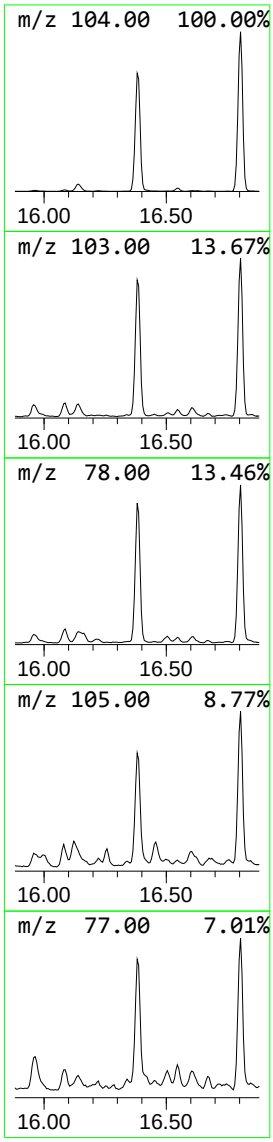
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 37 Benzene, 1,1'-(1,2-cyclobut... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.381	23.38 ng/ul	5180970	Phenanthrene-d10	17.204

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	93
2	[2.2]Paracyclophane	208	C16H16	001633-22-3	90
3	[2.2]Paracyclophane	208	C16H16	001633-22-3	87
4	Cyclobutane, 1,2-diphenyl-	208	C16H16	003018-21-1	78
5	5,5-Dimethyl-4-phenyldihydrofura...	190	C12H14O2	013133-96-5	78





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Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

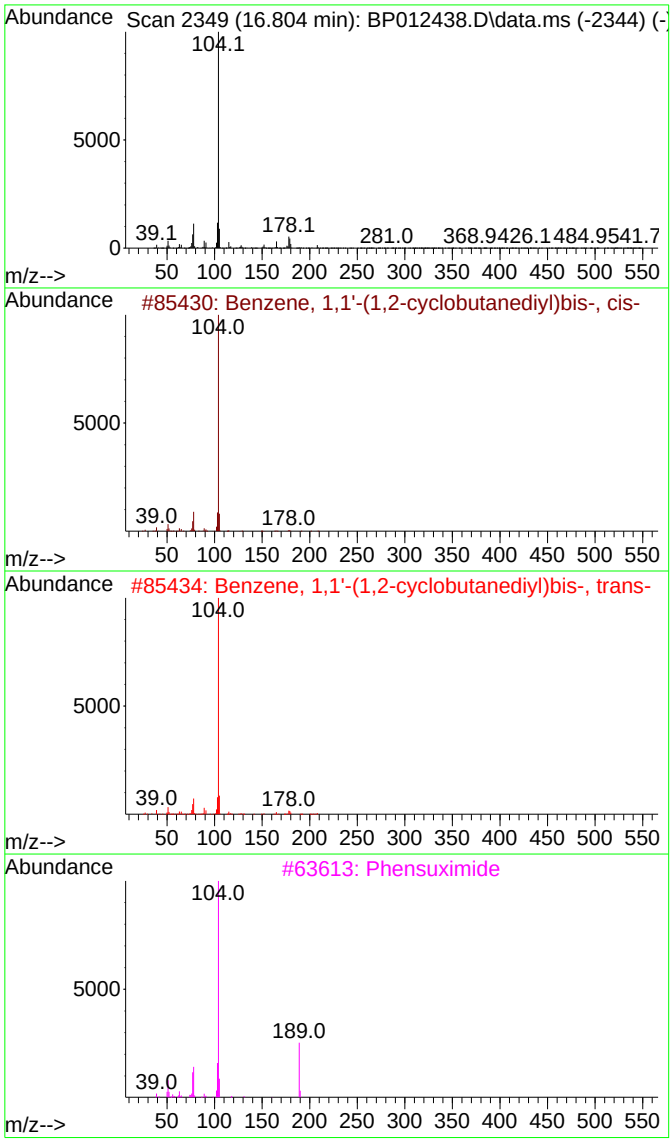
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

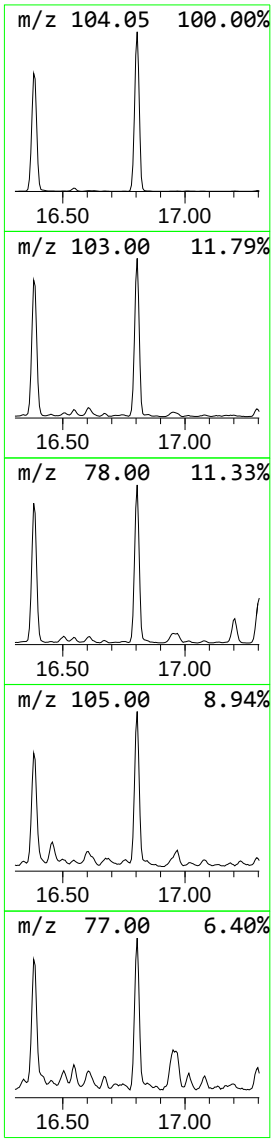
****

Peak Number 41 Phensuximide Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.804	26.79 ng/ul	5936670	Phenanthrene-d10	17.204

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	86
2	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	72
3	Phensuximide	189	C11H11NO2	000086-34-0	64
4	[2.2]Paracyclophane	208	C16H16	001633-22-3	64
5	[2.2]Paracyclophane	208	C16H16	001633-22-3	64





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Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

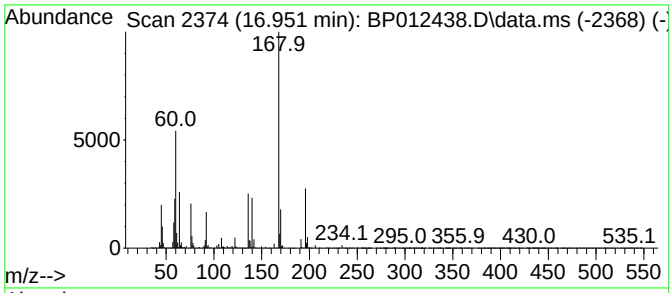
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

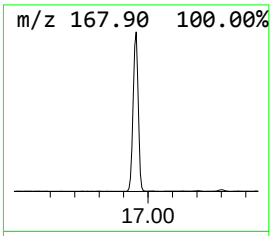
****

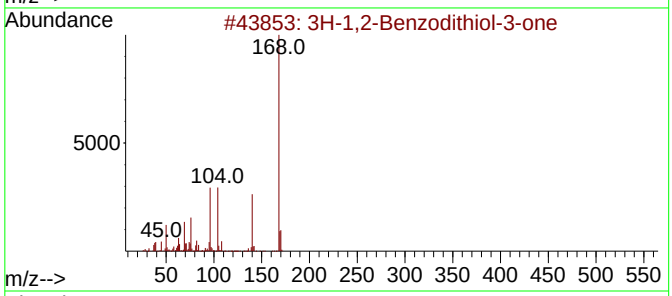
Peak Number 42 unknown-02 Concentration Rank 21

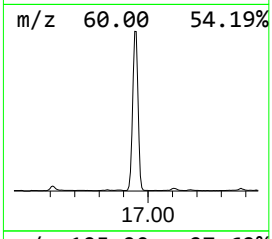
R.T.	EstConc	Area	Relative to ISTD	R.T.
16.951	22.50 ng/ul	4987270	Phenanthrene-d10	17.204

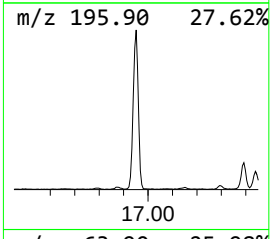
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3H-1,2-Benzodithiol-3-one	168	C7H4OS2	001677-27-6	38
2	(.eta.5-cyclopentadienyl)dicarbo...	224	C7H5O2Rh	012192-97-1	28
3	Rhodium, (.eta.5-2,4-cyclopentad...	224	C9H13Rh	012211-95-9	16
4	p-Nitrobenzoic acid, n-propyl ester	209	C10H11NO4	000094-22-4	12
5	Thiazole, 4,5-dihydro-2-methyl-	101	C4H7NS	002346-00-1	11

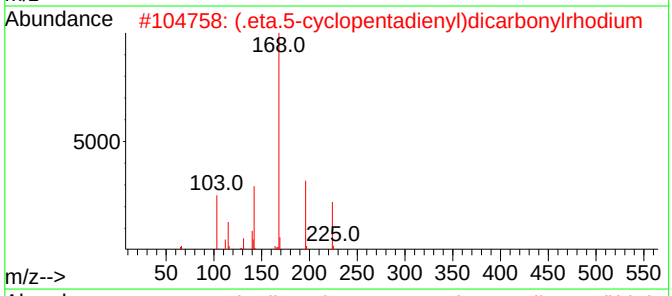


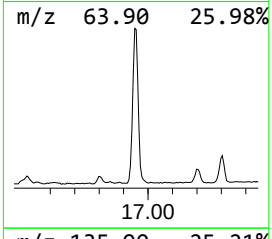


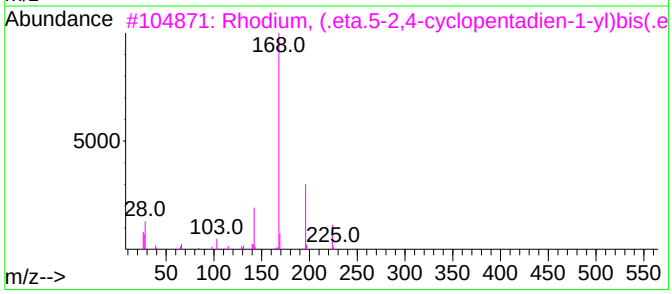


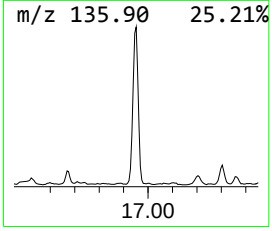












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

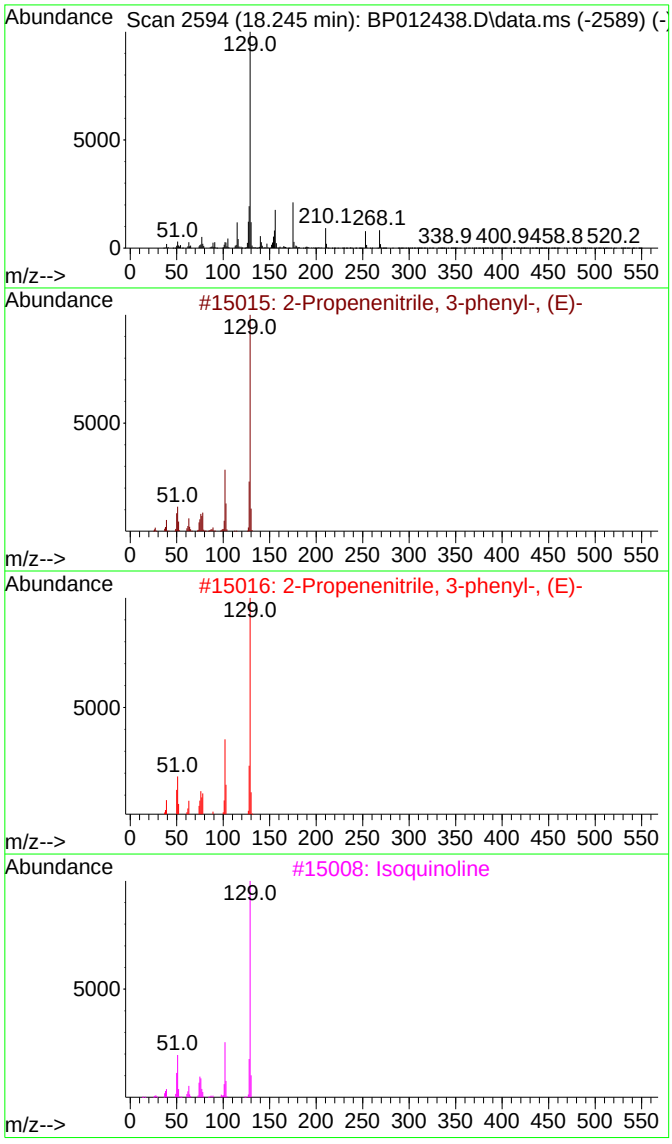
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

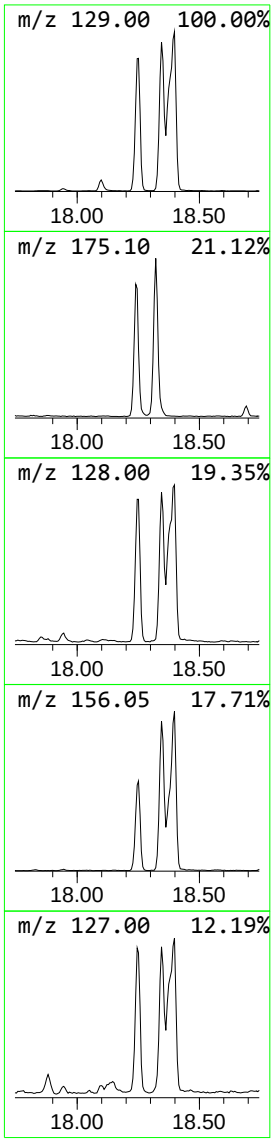
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 47 2-Propenenitrile, 3-phenyl-... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.245	18.37 ng/ul	4070590	Phenanthrene-d10	17.204

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenenitrile, 3-phenyl-, (E)-	129	C9H7N	001885-38-7	50
2	2-Propenenitrile, 3-phenyl-, (E)-	129	C9H7N	001885-38-7	50
3	Isoquinoline	129	C9H7N	000119-65-3	40
4	2-Chloro-1-(2-methyl-allyl)-1,2,...	220	C14H17Cl	1000185-82-6	38
5	Isoquinoline	129	C9H7N	000119-65-3	38





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Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

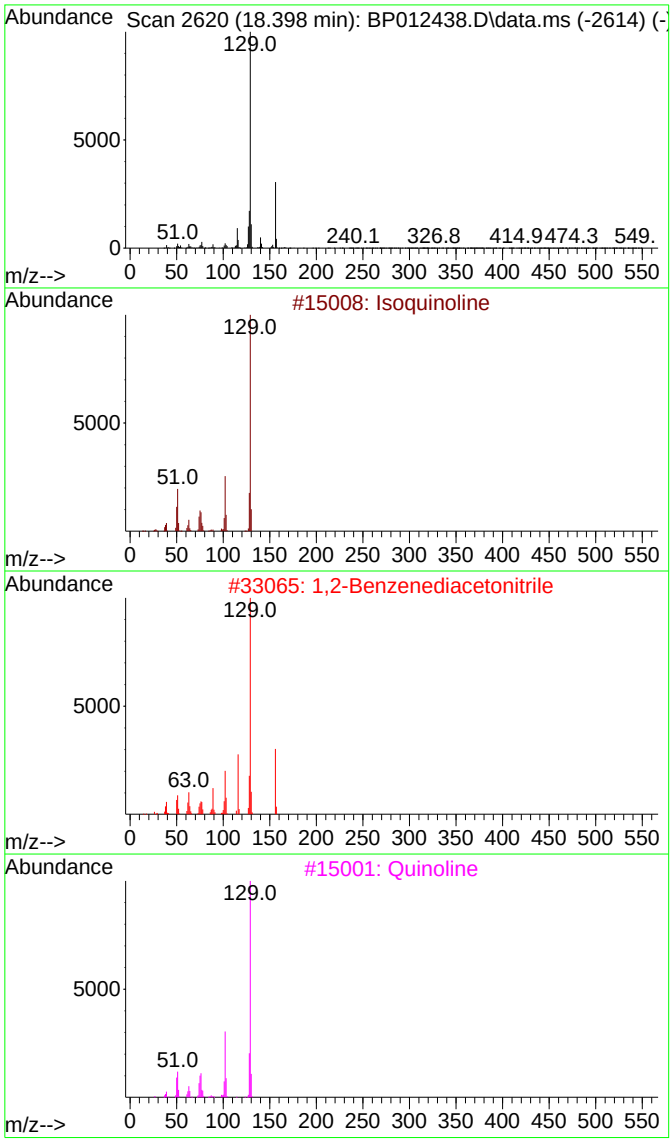
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

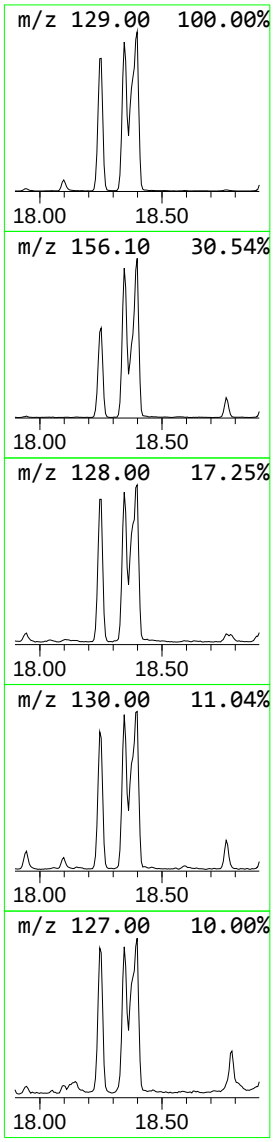
****

Peak Number 50 Isoquinoline Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.398	21.38 ng/ul	4738200	Phenanthrene-d10	17.204

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isoquinoline	129	C9H7N	000119-65-3	52
2	1,2-Benzenediacetonitrile	156	C10H8N2	000613-73-0	50
3	Quinoline	129	C9H7N	000091-22-5	49
4	Isoquinoline	129	C9H7N	000119-65-3	47
5	Isoquinoline	129	C9H7N	000119-65-3	47





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Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

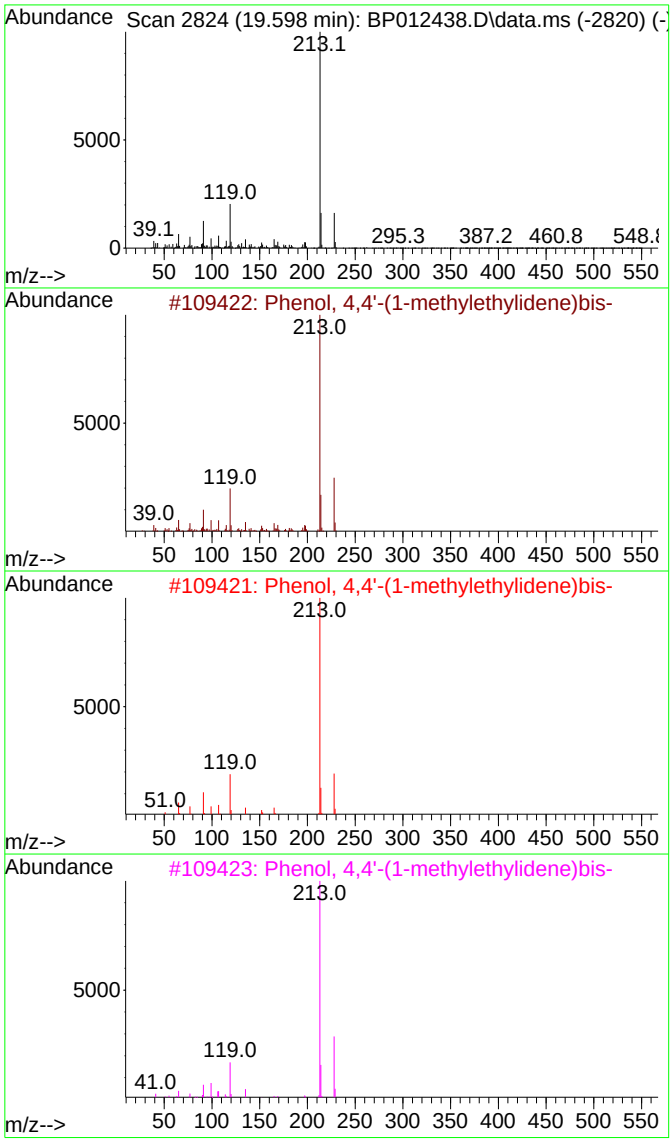
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

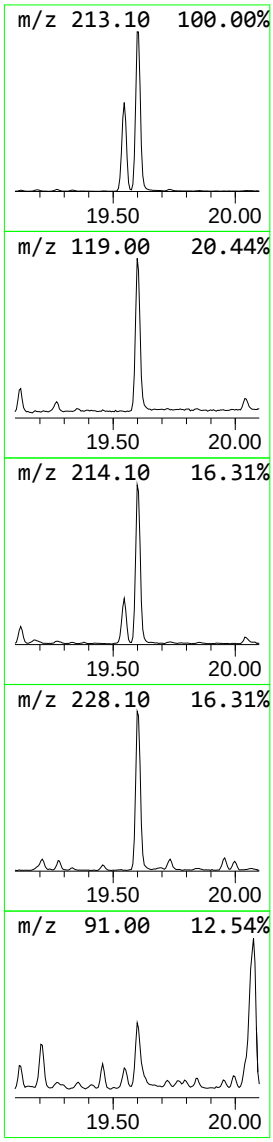
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 57 Phenol, 4,4'-(1-methylethyl... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.598	14.81 ng/ul	3888660	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	96
2	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	95
3	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	90
4	3,4'-Isopropylidenediphenol	228	C15H16O2	046765-25-7	78
5	4,4'-(1,3-Dimethylbutylidene)bis...	270	C18H22O2	006807-17-6	74





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Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

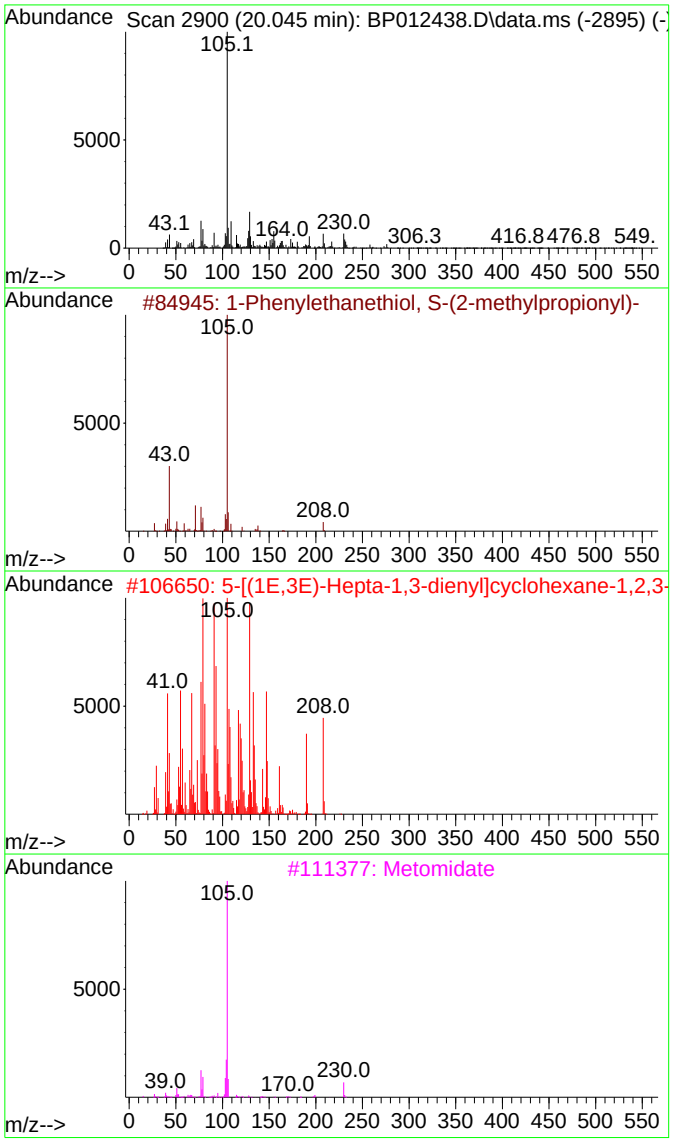
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

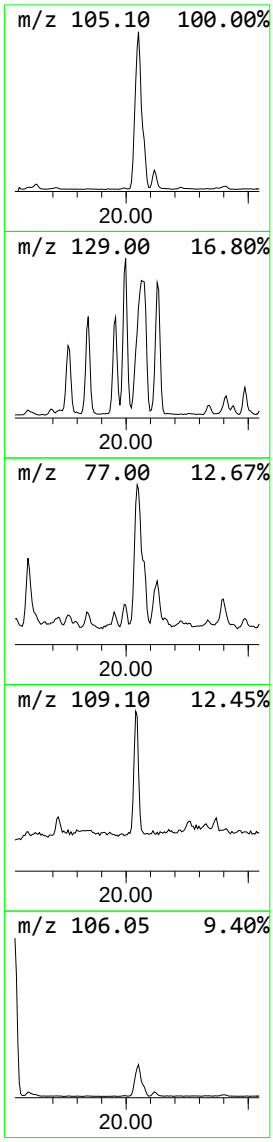
****

Peak Number 60 unknown-03 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.045	26.17 ng/ul	6868260	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenylethanethiol, S-(2-methyl...	208	C12H16OS	1227931-65-8	43
2	5-[(1E,3E)-Hepta-1,3-dienyl]cycl...	226	C13H22O3	2111875-27-3	42
3	Metomidate	230	C13H14N2O2	005377-20-8	41
4	Benzaldehyde, 3-methoxy-4-[(2-me...	256	C16H16O3	1000338-23-4	38
5	Benzene, 1,1'-(1,2-dimethyl-1,2-...	210	C16H18	005789-35-5	38





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Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

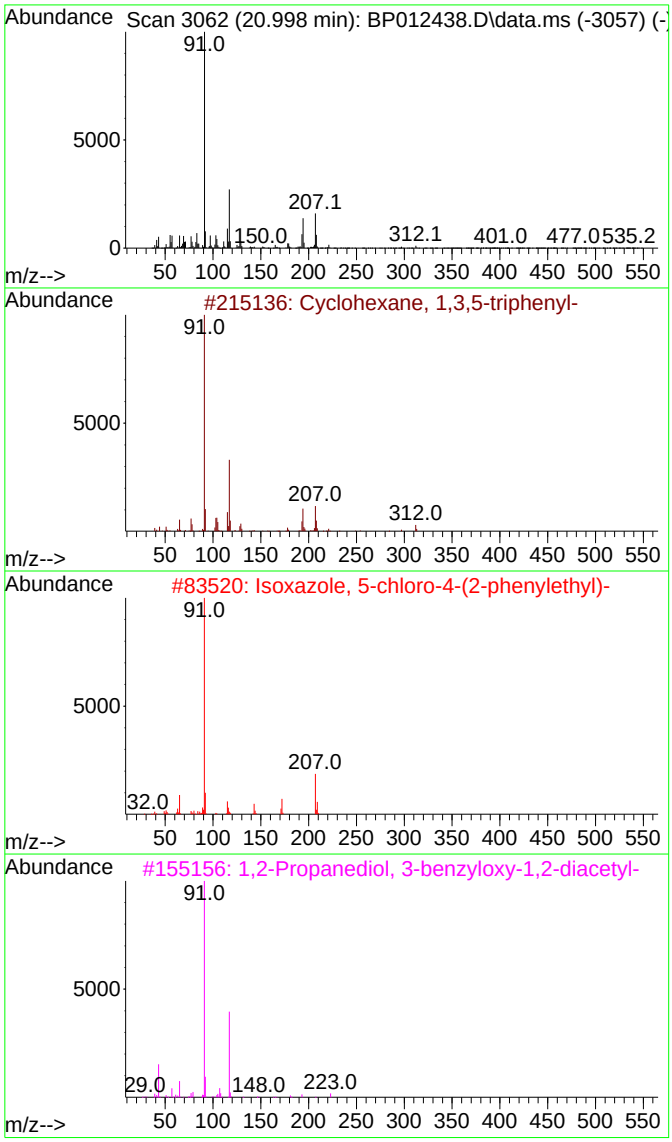
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

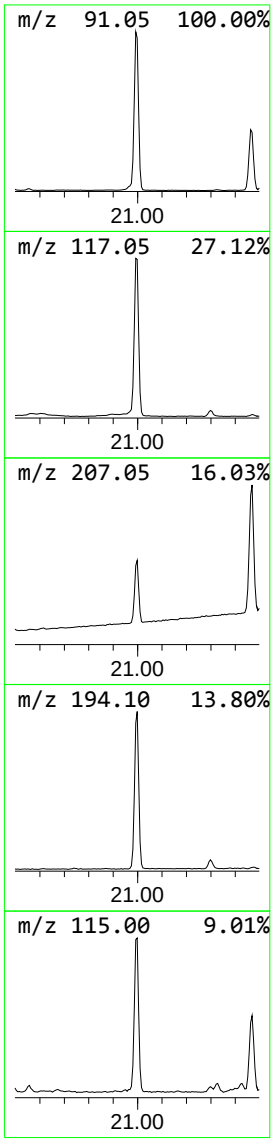
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 64 Cyclohexane, 1,3,5-triphenyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.998	34.10 ng/ul	8951540	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,3,5-triphenyl-	312	C24H24	028336-57-4	72
2	Isoxazole, 5-chloro-4-(2-phenyle...	207	C11H10ClNO	1000362-09-0	38
3	1,2-Propanediol, 3-benzyloxy-1,2...	266	C14H18O5	013754-10-4	25
4	Carbonic acid, monoamide, N-benz...	207	C12H17NO2	1000451-48-5	22
5	(R)-(1-Cyclopentylpropane-1,3-di...	264	C20H24	1402851-74-4	17





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Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

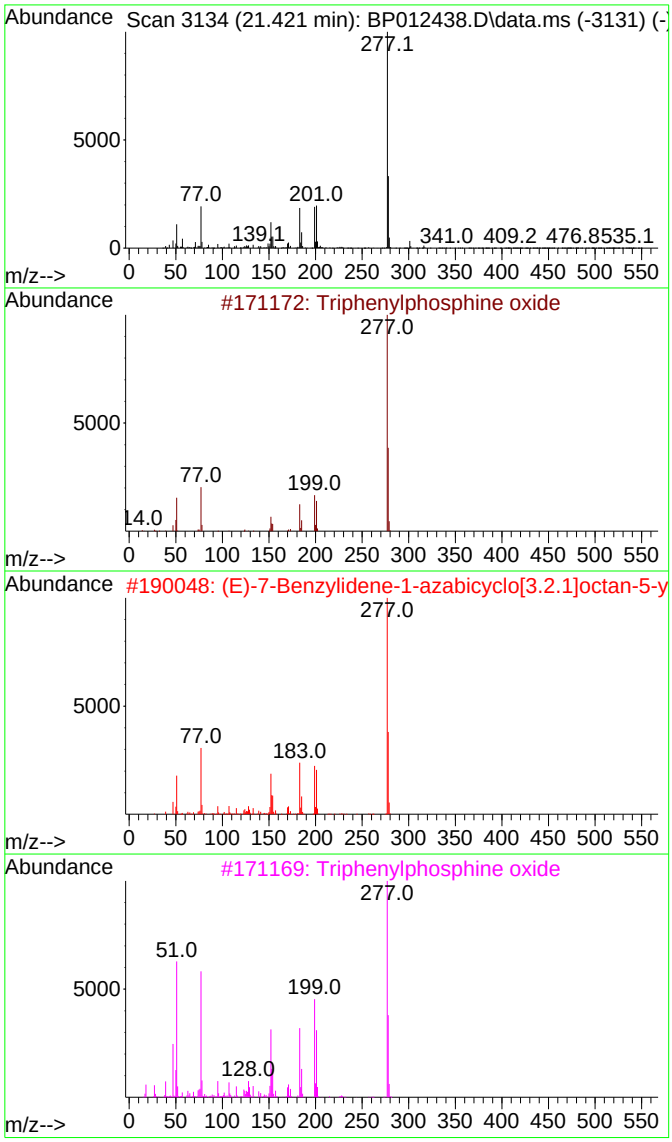
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

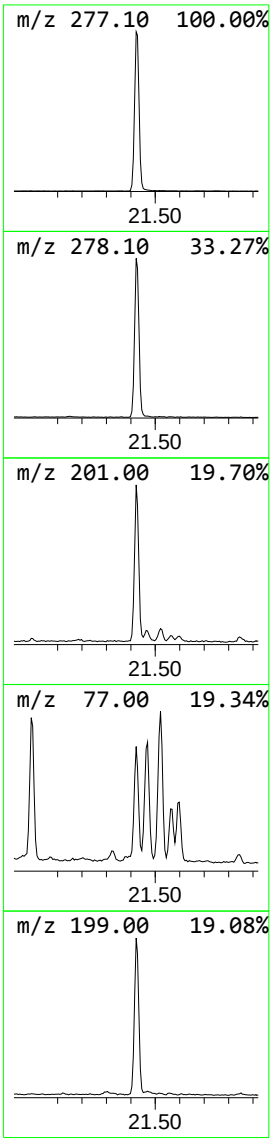
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 65 Triphenylphosphine oxide Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.422	7.53 ng/ul	1976050	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	93
2	(E)-7-Benzylidene-1-azabicyclo[3...	293	C15H19NO3S	1000483-83-4	91
3	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	87
4	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	60
5	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	49





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Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

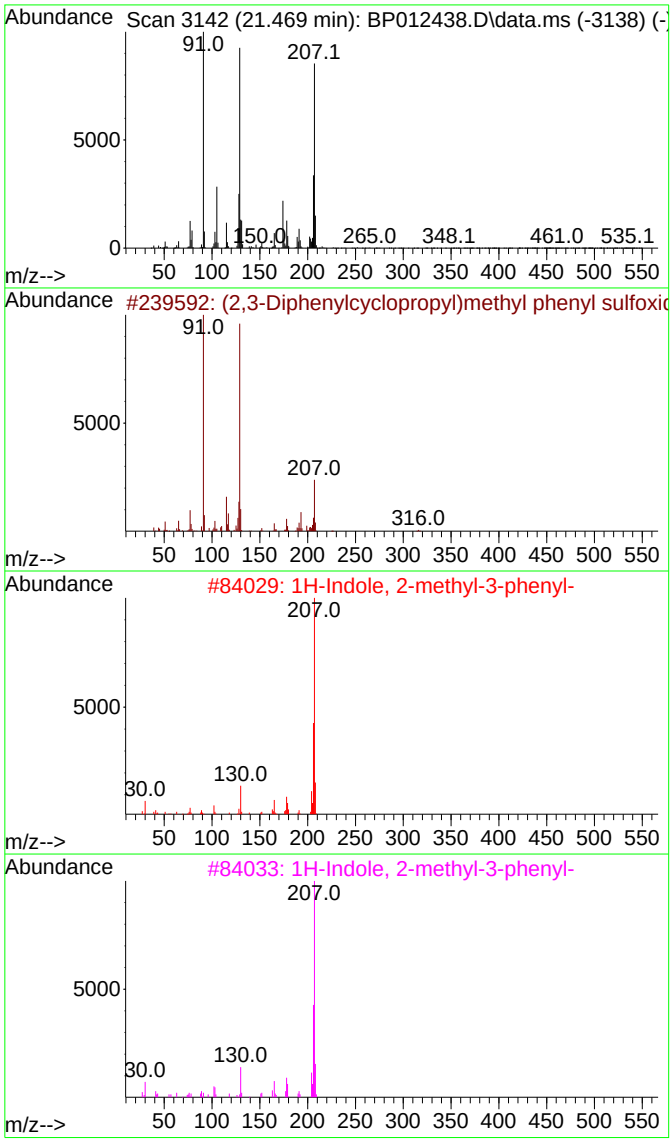
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

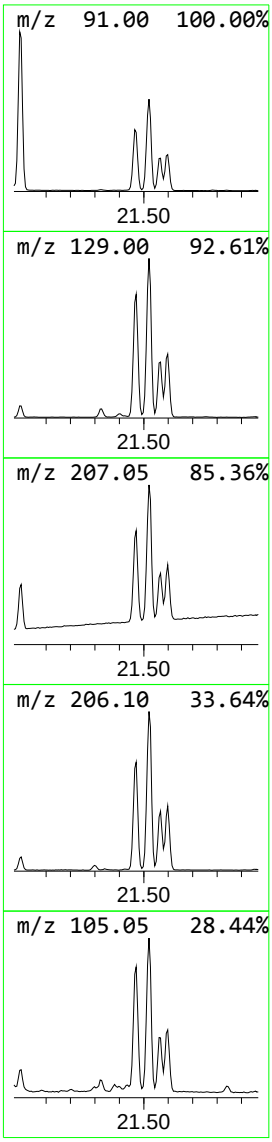
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 66 (2,3-Diphenylcyclopropyl)me... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.469	21.15 ng/ul	5552890	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	64
2	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	50
3	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	50
4	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	46
5	Cyclopentene-1-carboxylic acid, ...	332	C23H24O2	1000159-40-6	45





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Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

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Quant Title : SVOA CALIBRATION

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

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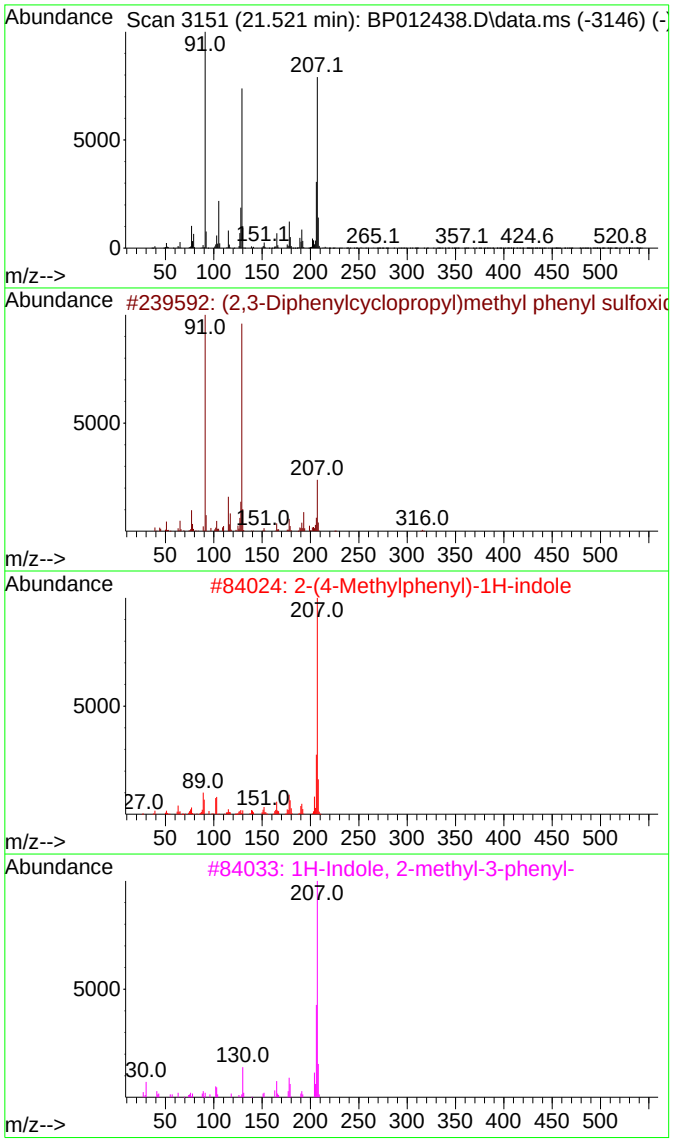
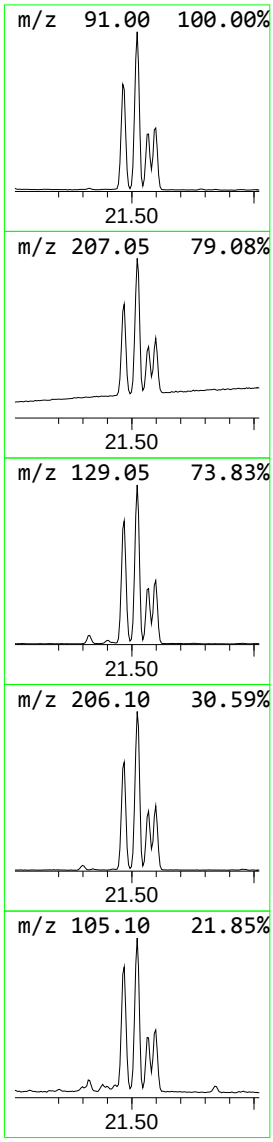
Peak Number 67 unknown-04 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.522	27.42 ng/ul	7197410	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	50
2	2-(4-Methylphenyl)-1H-indole	207	C15H13N	055577-25-8	46
3	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	43
4	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	43
5	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	41





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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

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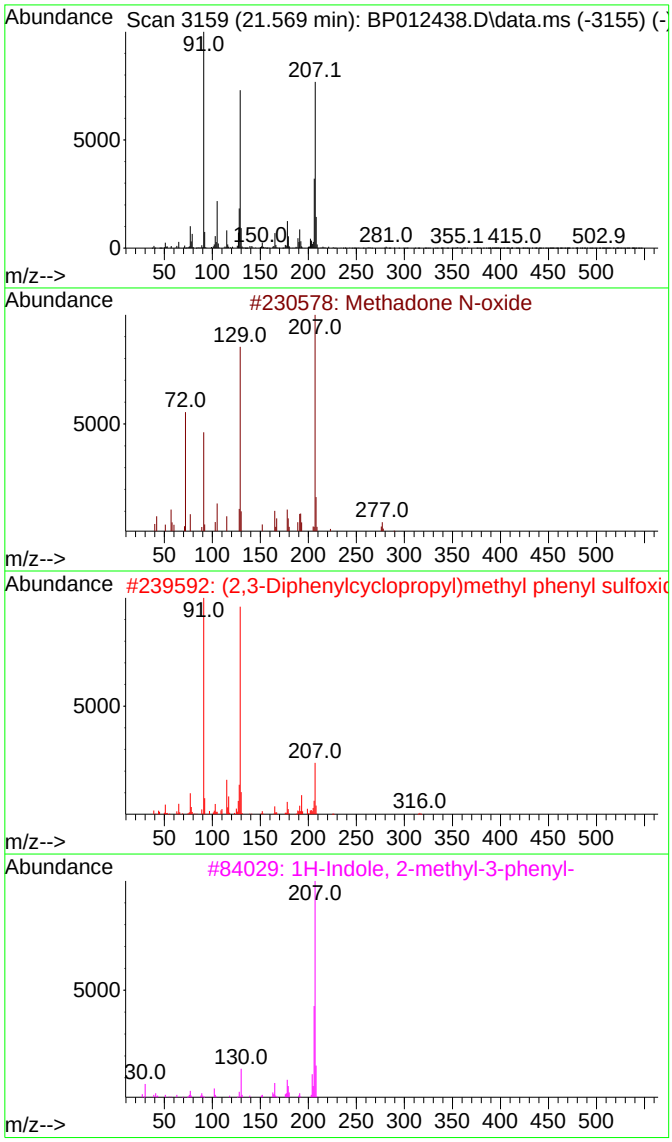
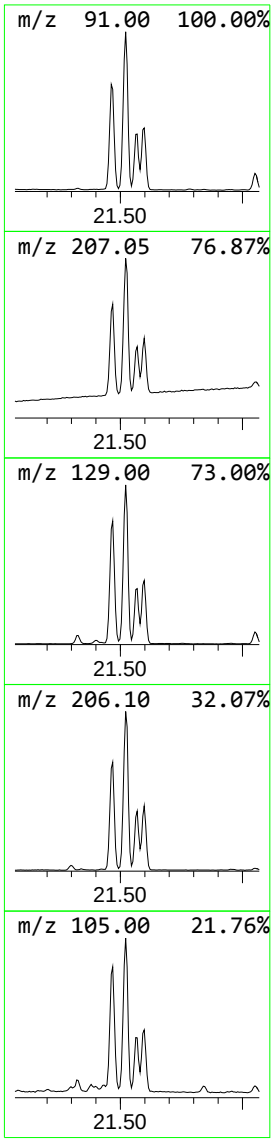
Peak Number 68 Methadone N-oxide Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.569	9.96 ng/ul	2615280	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methadone N-oxide	325	C21H27NO2	1000120-80-7	58
2	(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	50
3	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	49
4	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	46
5	5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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

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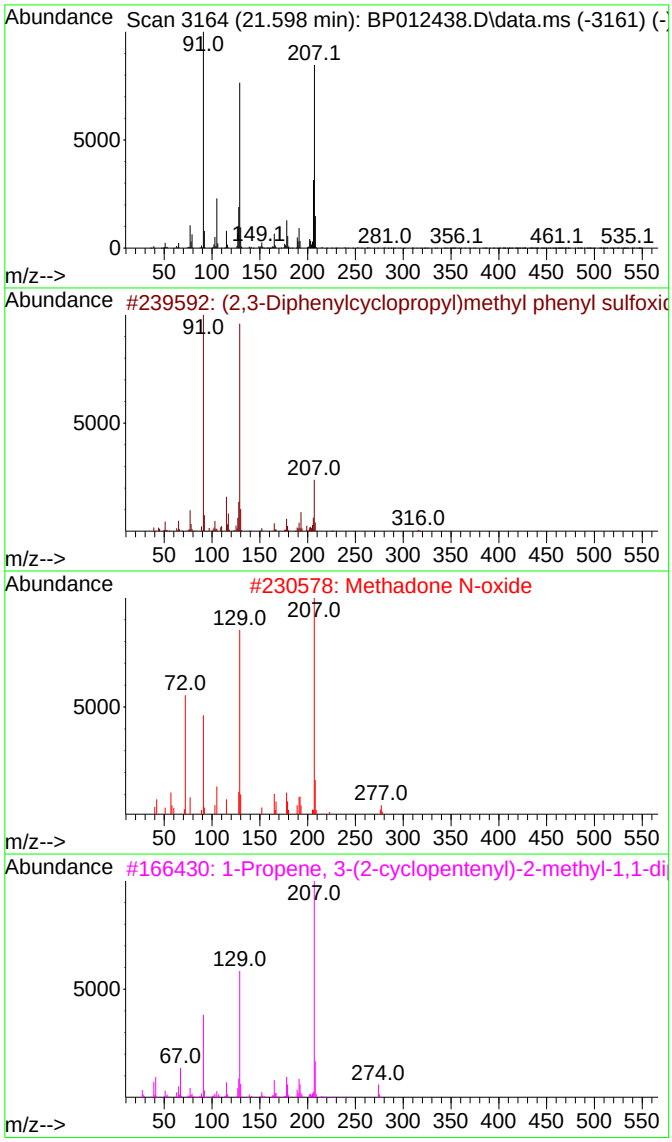
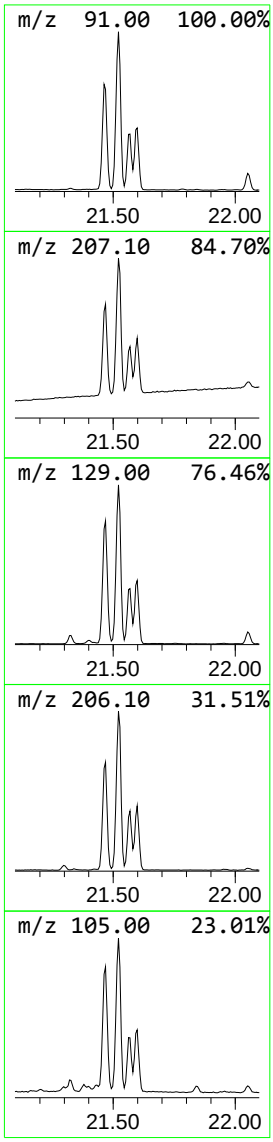
Peak Number 69 unknown-05 Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.598	10.35 ng/ul	2716740	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	72
2	Methadone N-oxide	325	C21H27NO2	1000120-80-7	64
3	1-Propene, 3-(2-cyclopentenyl)-2...	274	C21H22	1010154-23-3	47
4	1-benzylindole	207	C15H13N	1000334-31-3	38
5	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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

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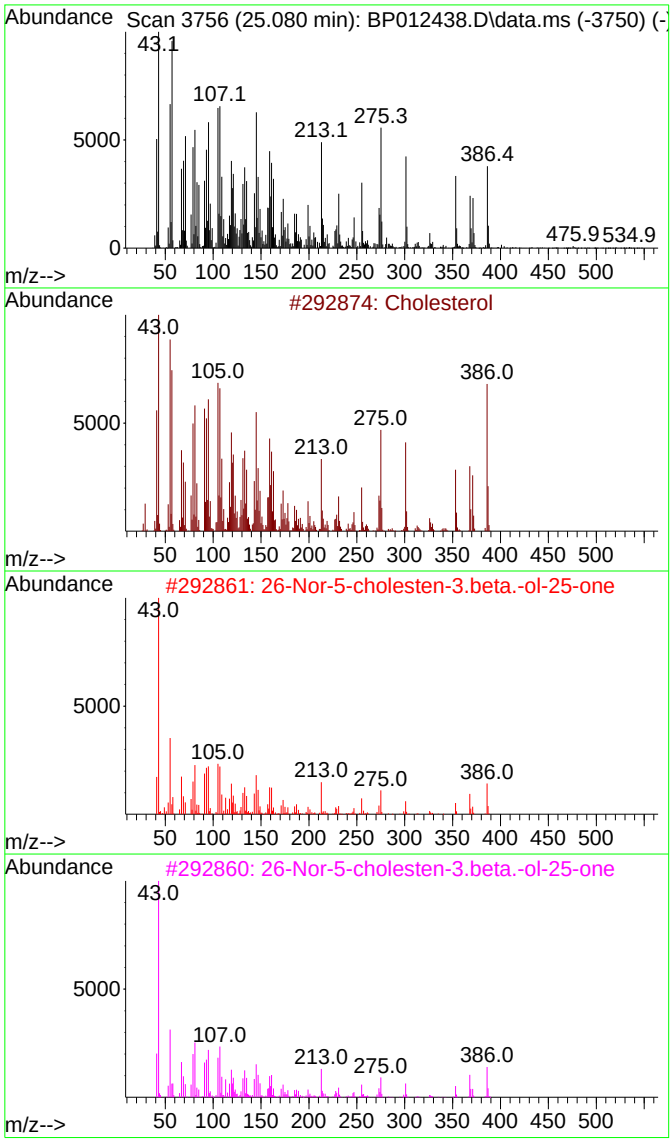
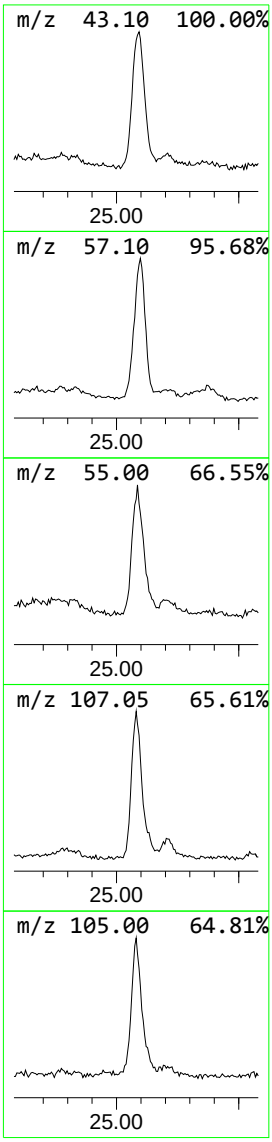
Peak Number 72 Cholesterol Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.080	14.21 ng/ul	3144530	Perylene-d12	23.686

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cholesterol	386	C27H46O	000057-88-5	99
2	26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	99
3	26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	96
4	Cholesterol	386	C27H46O	000057-88-5	95
5	Cholest-5-en-3-ol, (3.alpha.)-	386	C27H46O	000474-77-1	70





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012438.D
Acq On : 01 Nov 2022 21:19
Operator : CG/JU
Sample : N5300-03
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA80

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.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
Benzene, 1,3-di...	5.793	66.8	ng/ul	5460550	1	7.787	1636070	20.0
Butyl aldoxime,...	6.058	14.0	ng/ul	1147460	1	7.787	1636070	20.0
unknown-01	6.458	13.1	ng/ul	1072530	1	7.787	1636070	20.0
Benzene, propyl-	6.758	44.3	ng/ul	3625920	1	7.787	1636070	20.0
Benzene, 1,2,3-...	7.428	51.5	ng/ul	4216730	1	7.787	1636070	20.0
Benzene, (1-met...	7.540	15.9	ng/ul	1297830	1	7.787	1636070	20.0
2-Methyl-1,3-di...	7.675	20.9	ng/ul	1705370	1	7.787	1636070	20.0
Benzene, 1,2,4-...	7.899	26.3	ng/ul	2152600	1	7.787	1636070	20.0
Furan, tetrahyd...	7.981	15.1	ng/ul	1234690	1	7.787	1636070	20.0
Indane	8.157	46.6	ng/ul	3810940	1	7.787	1636070	20.0
Cyclohexanone, ...	8.222	65.7	ng/ul	5373840	1	7.787	1636070	20.0
Benzene, 1,3-di...	8.275	26.1	ng/ul	2136720	1	7.787	1636070	20.0
Cyclohexanol, 3...	8.416	93.4	ng/ul	7638630	1	7.787	1636070	20.0
o-Cymene	8.887	15.3	ng/ul	1247410	1	7.787	1636070	20.0
3-Octanol, 3,7-...	9.104	57.0	ng/ul	4665930	1	7.787	1636070	20.0
Phenol, p-tert-...	11.945	35.8	ng/ul	4141410	2	10.593	2311950	20.0
Benzene, 1,1'-(...	16.381	23.4	ng/ul	5180970	4	17.204	4432280	20.0
Phensuximide	16.804	26.8	ng/ul	5936670	4	17.204	4432280	20.0
unknown-02	16.951	22.5	ng/ul	4987270	4	17.204	4432280	20.0
2-Propenenitril...	18.245	18.4	ng/ul	4070590	4	17.204	4432280	20.0
Isoquinoline	18.398	21.4	ng/ul	4738200	4	17.204	4432280	20.0
Phenol, 4,4'-(1...	19.598	14.8	ng/ul	3888660	5	21.298	5249950	20.0
unknown-03	20.045	26.2	ng/ul	6868260	5	21.298	5249950	20.0
Cyclohexane, 1,...	20.998	34.1	ng/ul	8951540	5	21.298	5249950	20.0
Triphenylphosph...	21.422	7.5	ng/ul	1976050	5	21.298	5249950	20.0
(2,3-Diphenylcy...	21.469	21.1	ng/ul	5552890	5	21.298	5249950	20.0
unknown-04	21.522	27.4	ng/ul	7197410	5	21.298	5249950	20.0
Methadone N-oxide	21.569	10.0	ng/ul	2615280	5	21.298	5249950	20.0
unknown-05	21.598	10.4	ng/ul	2716740	5	21.298	5249950	20.0
Cholesterol	25.080	14.2	ng/ul	3144530	6	23.686	4425410	20.0