LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013747.D
Acq On : 03 Feb 2023 18:57
Operator : CG/JU
Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A44N2

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-BP020123.M
Title : SVOA CALIBRATION

Signal : TIC: BP013747.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.446	38	44	68	rVB	5765776	10406466	25.41%	5.609%
2	3.917	121	124	139	rBV	132725	247391	0.60%	0.133%
3	4.246	175	180	191	rBV	86031	140678	0.34%	0.076%
4	5.411	370	378	394	rBV	13747090	23638552	57.72%	12.741%
5	5.605	404	411	425	rBV2	80976	152385	0.37%	0.082%
6	6.116	492	498	505	rBV	39243	65078	0.16%	0.035%
7	6.469	552	558	564	rVB	27029	44273	0.11%	0.024%
8	6.775	605	610	616	rBV2	29675	58107	0.14%	0.031%
9	7.099	658	665	671	rBV	58697	96616	0.24%	0.052%
10	7.275	689	695	705	rBV	147926	245370	0.60%	0.132%
11	7.452	718	725	731	rBV	739047	1187053	2.90%	0.640%
12	7.511	731	735	740	rVB	37280	57326	0.14%	0.031%
13	7.658	753	760	770	rVB	605369	983116	2.40%	0.530%
14	7.769	775	779	786	rVB2	26945	46867	0.11%	0.025%
15	8.022	815	822	830	rVB	1009355	1637287	4.00%	0.882%
16	8.134	830	841	842	rBV	725159	1006043	2.46%	0.542%
17	8.169	842	847	856	rVV	6846430	11246838	27.46%	6.062%
18	8.246	856	860	865	rVB	29720	45311	0.11%	0.024%
19	8.505	889	904	912	rBV	21471872	40952903	100.00%	22.073%
20	8.834	948	960	968	rBV	444518	733873	1.79%	0.396%
21	9.305	1032	1040	1043	rBV	736074	1270401	3.10%	0.685%
22	9.346	1043	1047	1061	rVB	1894455	3165834	7.73%	1.706%
23	9.705	1104	1108	1114	rVB2	27368	42903	0.10%	0.023%
24	9.975	1147	1154	1158	rBV	151210	256414	0.63%	0.138%
25	10.034	1158	1164	1179	rVB	601702	1035417	2.53%	0.558%
26	10.240	1193	1199	1206	rBV	176632	278450	0.68%	0.150%
27	10.569	1247	1255	1263	rBV	1002935	1735021	4.24%	0.935%
28	10.640	1263	1267	1273	rVV3	24511	49569	0.12%	0.027%
29	10.822	1290	1298	1308	rBV	2733705	4571600	11.16%	2.464%
30	10.957	1314	1321	1328	rBV	989353	1732201	4.23%	0.934%
31	11.010	1328	1330	1336	rVV	36338	47471	0.12%	0.026%
32	11.093	1336	1344	1362	rVB	666753	1225618	2.99%	0.661%
33	11.346	1377	1387	1394	rBV	529393	871292	2.13%	0.470%
34	12.469	1569	1578	1584	rBV	305649	500593	1.22%	0.270%
35	12.975	1652	1664	1680	rBV	1343158	2084327	5.09%	1.123%
36	13.098	1680	1685	1692	rVV	31915	60347	0.15%	0.033%

LSC Area Percent Report

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Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

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-BP020123.M
Title : SVOA CALIBRATION

37	13.198	1695	1702	1708	rVV8	14916	43045	0.11%	0.023%
38	13.275	1711	1715	1724	rVV4	28629	61347	0.15%	0.033%
39	13.604	1768	1771	1775	rVV2	42303	79252	0.19%	0.043%
40	13.651	1775	1779	1784	rVB	95779	145314	0.35%	0.078%
41	13.816	1804	1807	1813	rVB	31850	44464	0.11%	0.024%
42	13.910	1819	1823	1829	rVB2	30318	49564	0.12%	0.027%
43	14.169	1860	1867	1882	rBV	2192352	3114829	7.61%	1.679%
44	14.463	1910	1917	1927	rVB	2544108	3636714	8.88%	1.960%
45	14.769	1962	1969	1975	rBV	1778834	2671674	6.52%	1.440%
46	14.951	1994	2000	2012	rBV	692507	1028450	2.51%	0.554%
47	15.257	2046	2052	2062	rVB2	115358	197737	0.48%	0.107%
48	15.757	2130	2137	2144	rVV	3746567	5246774	12.81%	2.828%
49	15.869	2150	2156	2167	rVB	734899	1047847	2.56%	0.565%
50	15.992	2172	2177	2184	rVB6	41424	84260	0.21%	0.045%
51	16.122	2192	2199	2205	rBV	2037288	2839670	6.93%	1.531%
52	16.257	2216	2222	2231	rBV	197372	303534	0.74%	0.164%
53	16.345	2231	2237	2248	rVV	787196	1131486	2.76%	0.610%
54	16.439	2248	2253	2258	rVV2	45883	100966	0.25%	0.054%
55	16.687	2290	2295	2306	rVB	120417	185563	0.45%	0.100%
56	16.810	2312	2316	2321	rBV	32513	48523	0.12%	0.026%
57	16.887	2325	2329	2337	rVB	83012	120478	0.29%	0.065%
58	17.269	2389	2394	2399	rBV	231155	341698	0.83%	0.184%
59	17.310	2399	2401	2409	rVV8	38692	77528	0.19%	0.042%
60	17.392	2410	2415	2423	rVB	303555	438919	1.07%	0.237%
61	17.522	2430	2437	2443	rBV	2612167	3749929	9.16%	2.021%
62	17.622	2447	2454	2460	rVV	4134385	6080319	14.85%	3.277%
63	17.692	2464	2466	2471	rVB2	43188	50645	0.12%	0.027%
64	17.798	2480	2484	2487	rVV	114765	167202	0.41%	0.090%
65	17.834	2487	2490	2494	rVB	73599	90327	0.22%	0.049%
66	18.375	2577	2582	2592	rBV	425626	614606	1.50%	0.331%
67	18.463	2593	2597	2604	rVB2	57959	117875	0.29%	0.064%
68	18.698	2632	2637	2643	rVB	270621	372816	0.91%	0.201%
69	19.110	2703	2707	2711	rBV	128287	168496	0.41%	0.091%
70	19.198	2718	2722	2725	rBV3	92294	130434	0.32%	0.070%
71	19.345	2744	2747	2751	rBV2	40412	48104	0.12%	0.026%
72	19.410	2754	2758	2761	rBV2	184327	257386	0.63%	0.139%
73	19.451	2761	2765	2768	rBV	852088	1020176	2.49%	0.550%
74	19.616	2786	2793	2803	rVB	358386	621410	1.52%	0.335%
75	19.833	2825	2830	2837	rBV2	7722815	10342118	25.25%	5.574%
76	19.898	2838	2841	2853	rVB3	161649	350182	0.86%	0.189%

LSC Area Percent Report

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Acq On : 03 Feb 2023 18:57
Operator : CG/JU
Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

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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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Title : SVOA CALIBRATION

77	20.104	2873	2876	2882	rVB	65496	90487	0.22%	0.049%
78	20.169	2884	2887	2891	rBV2	63422	80065	0.20%	0.043%
79	20.222	2893	2896	2902	rVB3	88498	121974	0.30%	0.066%
80	20.751	2982	2986	2992	rBV	293177	415234	1.01%	0.224%
81	20.898	3008	3011	3015	rBV2	104836	123728	0.30%	0.067%
82	21.275	3070	3075	3086	rBV3	3078573	4980415	12.16%	2.684%
83	21.598	3125	3130	3138	rVB	3781768	4973497	12.14%	2.681%
84	22.898	3348	3351	3359	rVB	371258	527836	1.29%	0.285%
85	23.998	3531	3538	3553	rVB2	4365082	9506484	23.21%	5.124%
86	24.168	3559	3567	3578	rVB2	2578914	5570660	13.60%	3.003%

Sum of corrected areas: 185531032

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013747.D
Acq On : 03 Feb 2023 18:57
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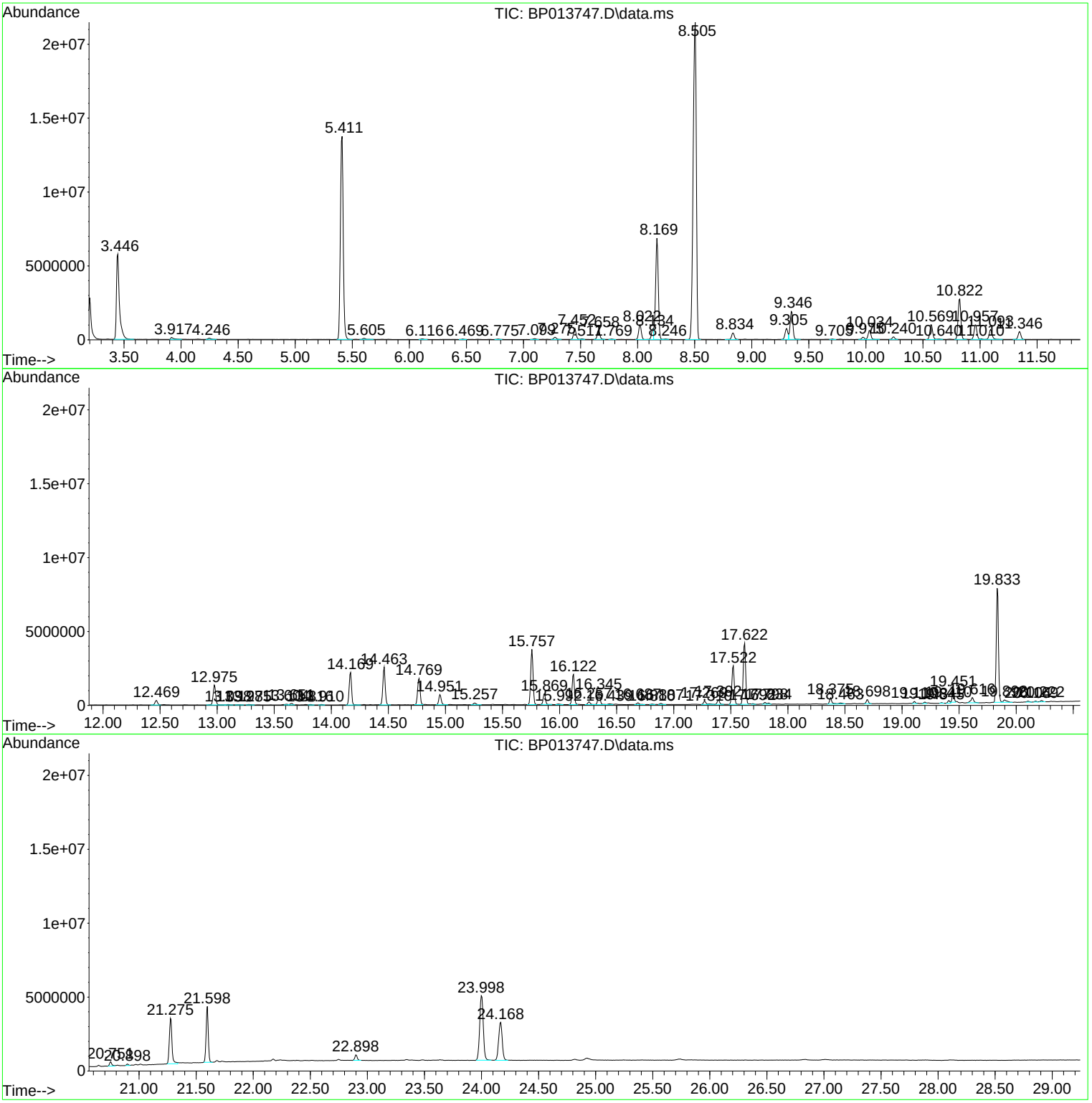
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013747.D
Acq On : 03 Feb 2023 18:57
Operator : CG/JU
Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

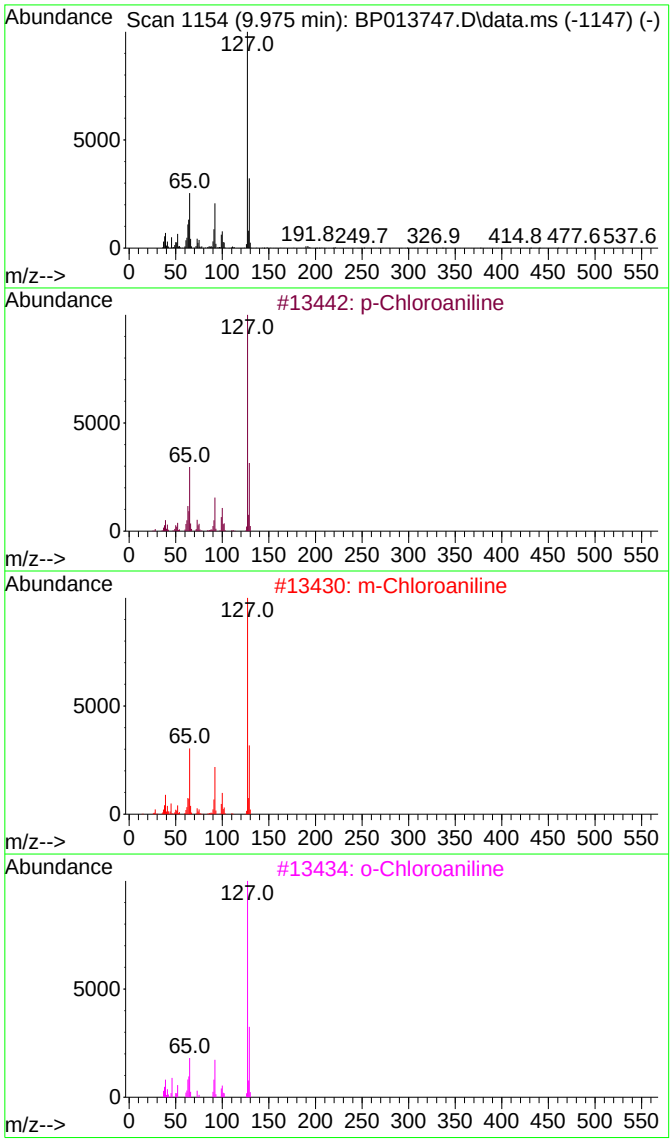
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

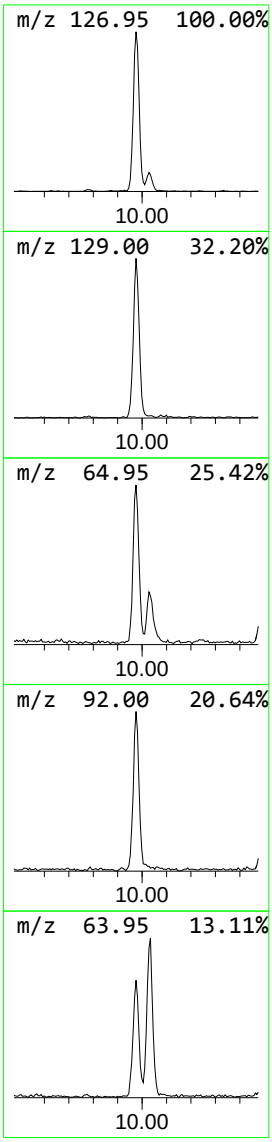
****

Peak Number 6 m-Chloroaniline Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.975	2.96 ng/ul	256414	Naphthalene-d8	10.957

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Chloroaniline	127	C6H6ClN	000106-47-8	97
2	m-Chloroaniline	127	C6H6ClN	000108-42-9	97
3	o-Chloroaniline	127	C6H6ClN	000095-51-2	97
4	o-Chloroaniline	127	C6H6ClN	000095-51-2	96
5	o-Chloroaniline	127	C6H6ClN	000095-51-2	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013747.D
Acq On : 03 Feb 2023 18:57
Operator : CG/JU
Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

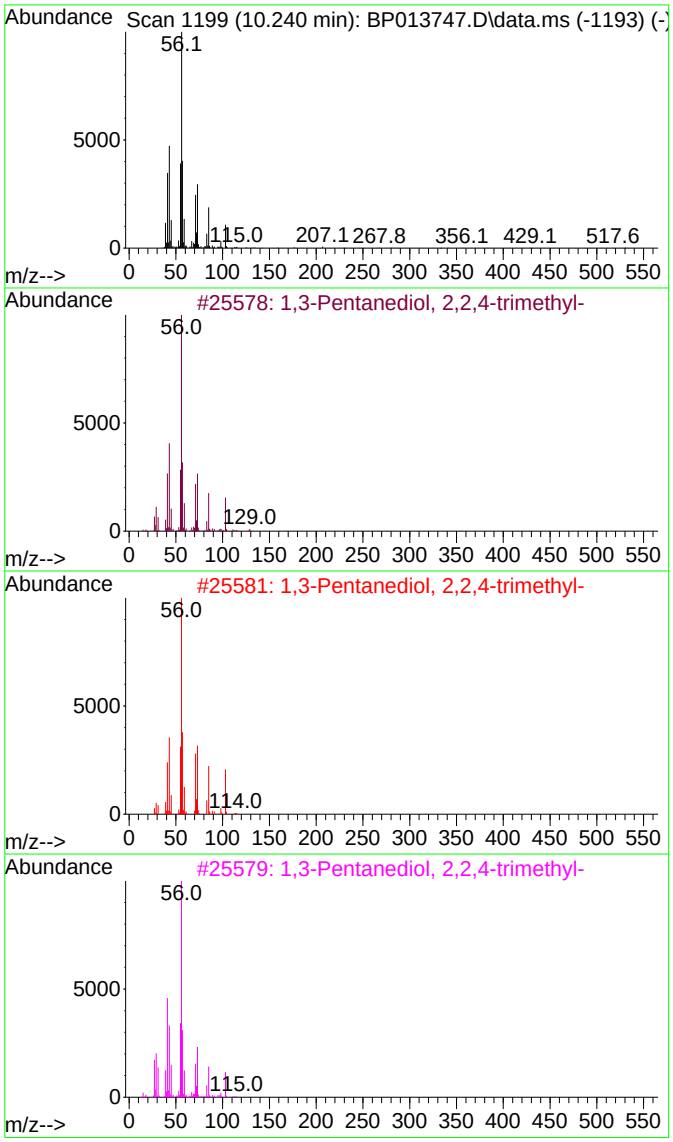
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

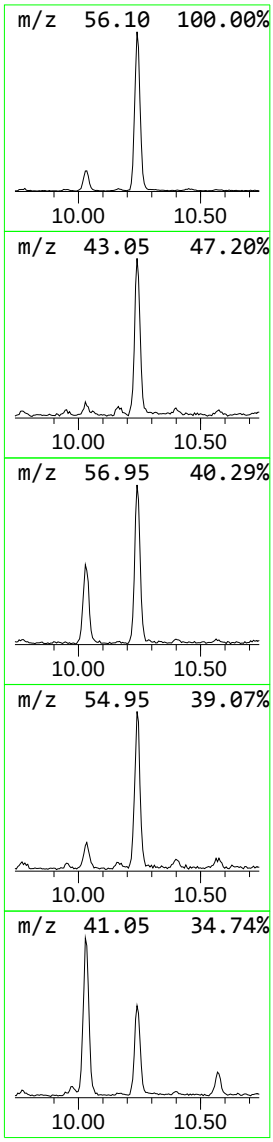
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1,3-Pentanediol, 2,2,4-trim... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.240	3.21 ng/ul	278450	Naphthalene-d8	10.957

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	86
2	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	64
3	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	59
4	2,3-Pentanediol, 2,4-dimethyl-	132	C7H16O2	066225-53-4	59
5	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
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Operator : CG/JU
Sample : 01381-13
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ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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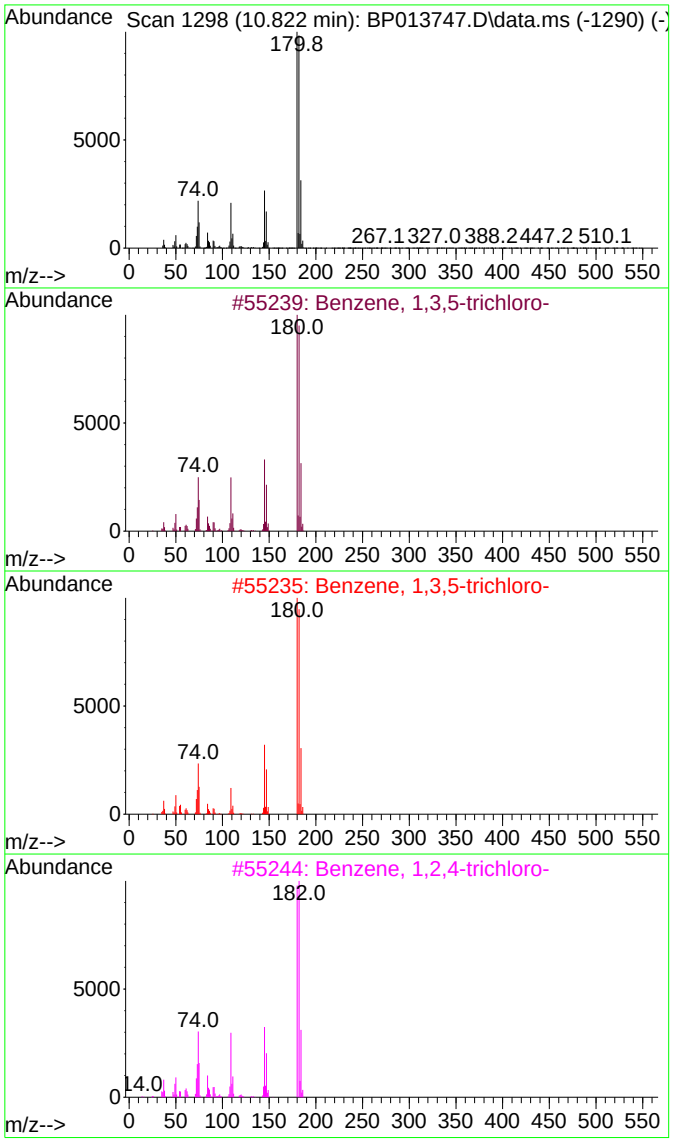
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

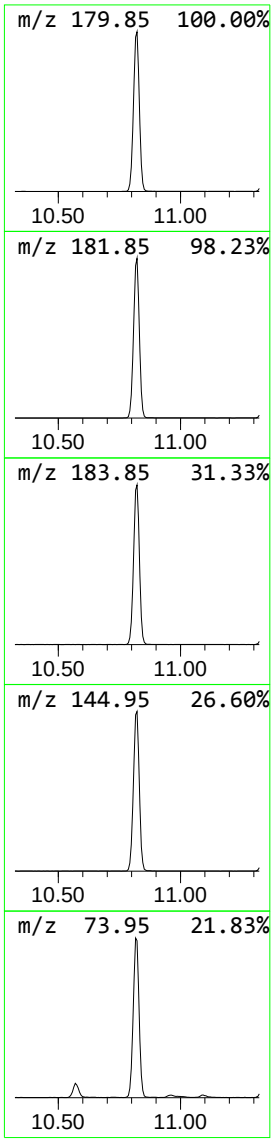
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,3,5-trichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.822	52.78 ng/ul	4571600	Naphthalene-d8	10.957

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
2	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3	Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
4	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97
5	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013747.D
Acq On : 03 Feb 2023 18:57
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Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

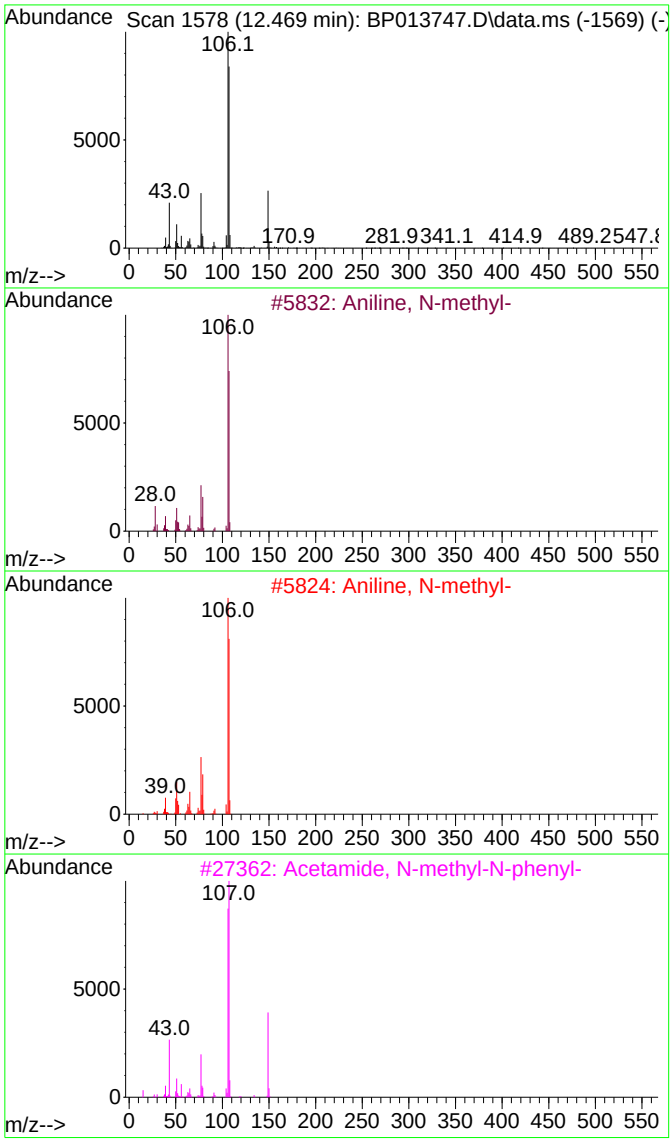
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

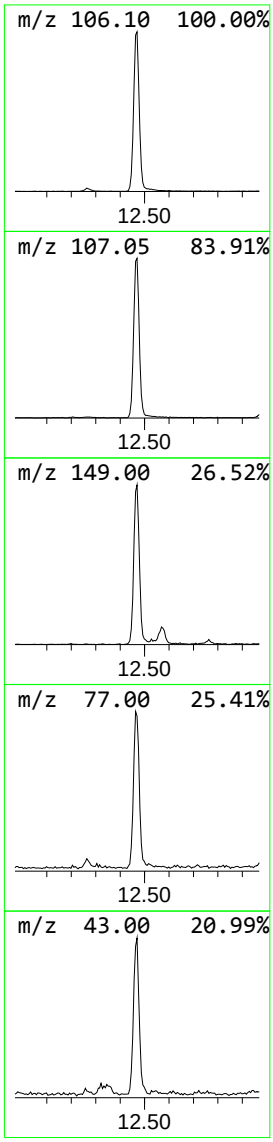
****

Peak Number 10 Aniline, N-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.469	5.78 ng/ul	500593	Naphthalene-d8	10.957

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Aniline, N-methyl-	107	C7H9N	000100-61-8	83
2	Aniline, N-methyl-	107	C7H9N	000100-61-8	83
3	Acetamide, N-methyl-N-phenyl-	149	C9H11NO	000579-10-2	81
4	Aniline, N-methyl-	107	C7H9N	000100-61-8	74
5	p-Aminotoluene	107	C7H9N	000106-49-0	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013747.D
Acq On : 03 Feb 2023 18:57
Operator : CG/JU
Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

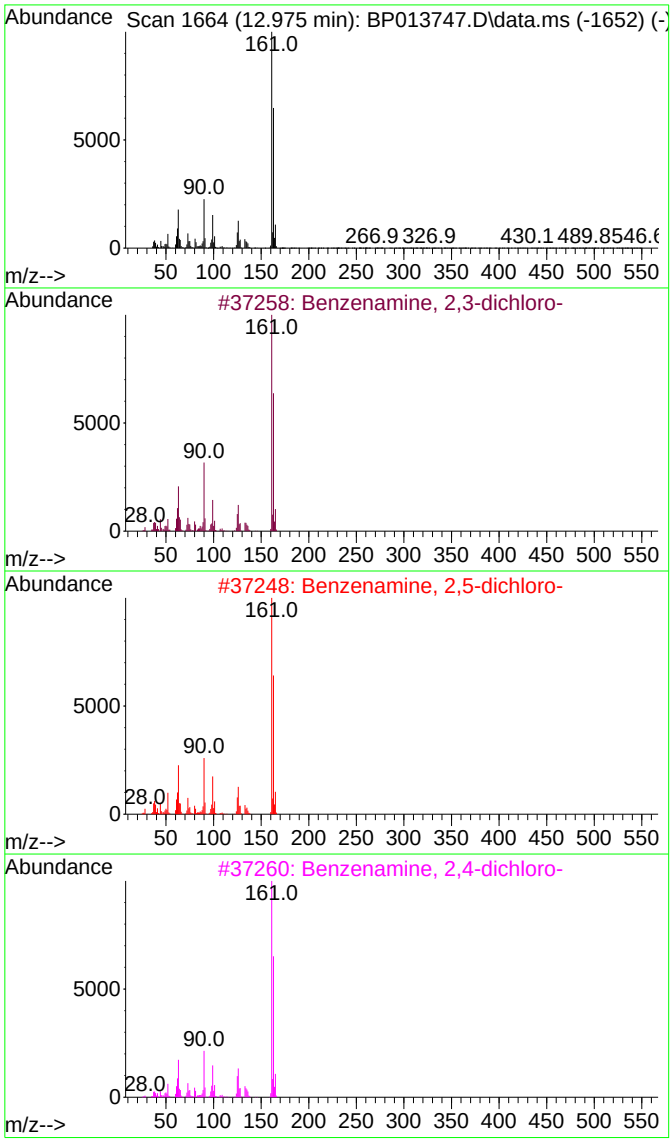
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

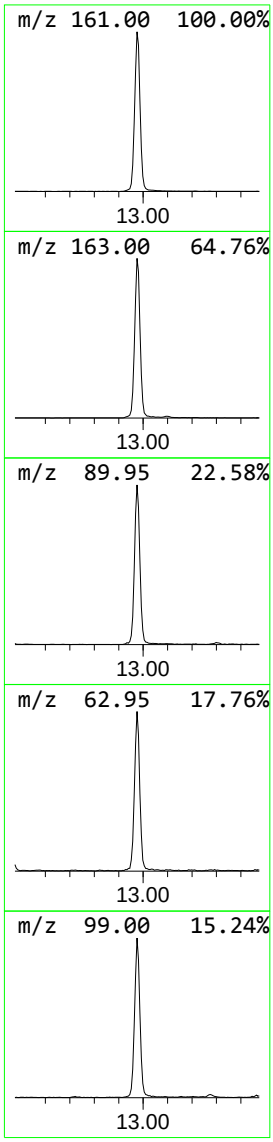
****

Peak Number 11 Benzenamine, 2,3-dichloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.975	15.60 ng/ul	2084330	Acenaphthene-d10	14.769

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 2,3-dichloro-	161	C6H5Cl2N	000608-27-5	98
2	Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	98
3	Benzenamine, 2,4-dichloro-	161	C6H5Cl2N	000554-00-7	97
4	Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	97
5	Benzenamine, 2,4-dichloro-	161	C6H5Cl2N	000554-00-7	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013747.D
Acq On : 03 Feb 2023 18:57
Operator : CG/JU
Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

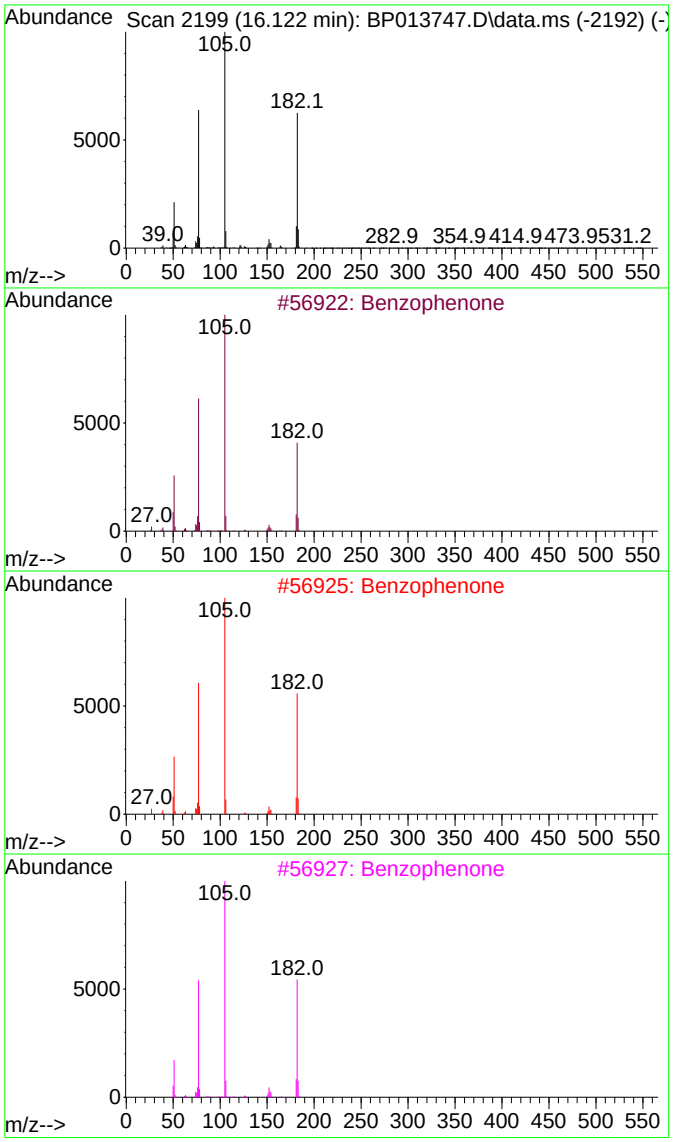
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

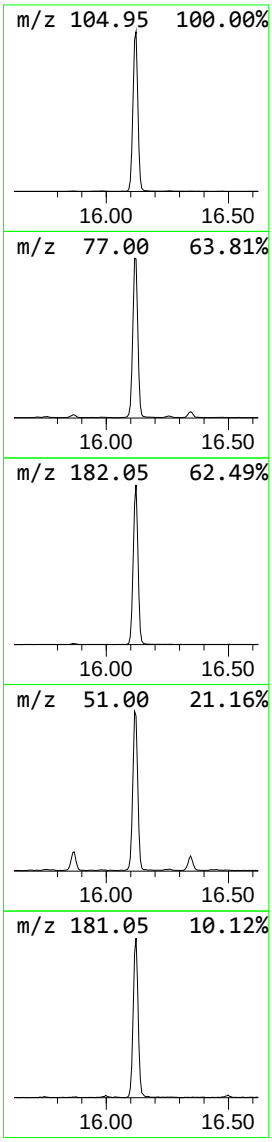
****

Peak Number 12 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.122	21.26 ng/ul	2839670	Acenaphthene-d10	14.769

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	97
2	Benzophenone	182	C13H10O	000119-61-9	96
3	Benzophenone	182	C13H10O	000119-61-9	96
4	Benzophenone	182	C13H10O	000119-61-9	95
5	Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	49





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Data File : BP013747.D
Acq On : 03 Feb 2023 18:57
Operator : CG/JU
Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

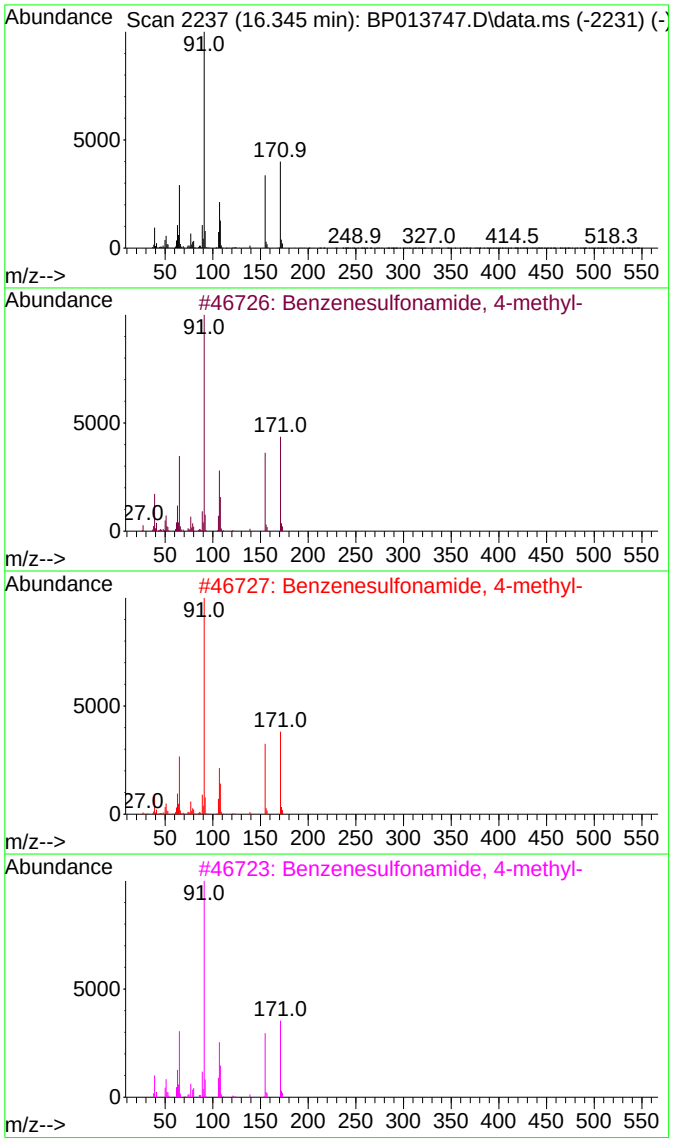
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

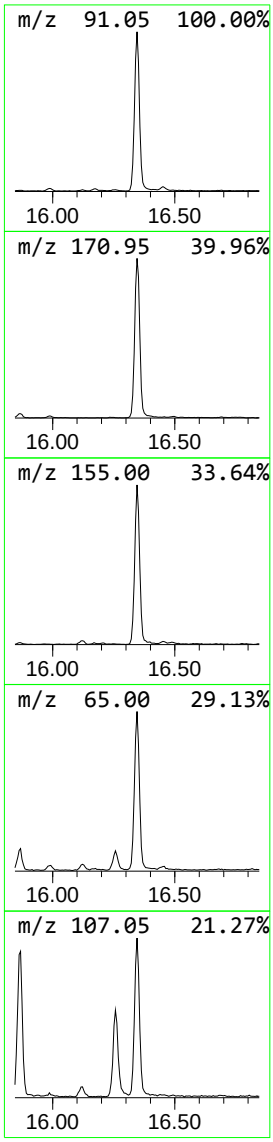
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzenesulfonamide, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.345	6.03 ng/ul	1131490	Phenanthrene-d10	17.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, 4-methyl-	171	C7H9NO2S	000070-55-3	98
2	Benzenesulfonamide, 4-methyl-	171	C7H9NO2S	000070-55-3	98
3	Benzenesulfonamide, 4-methyl-	171	C7H9NO2S	000070-55-3	98
4	Benzenesulfonamide, 4-methyl-	171	C7H9NO2S	000070-55-3	97
5	Benzenesulfonamide, 4-methyl-	171	C7H9NO2S	000070-55-3	94





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Acq On : 03 Feb 2023 18:57
Operator : CG/JU
Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

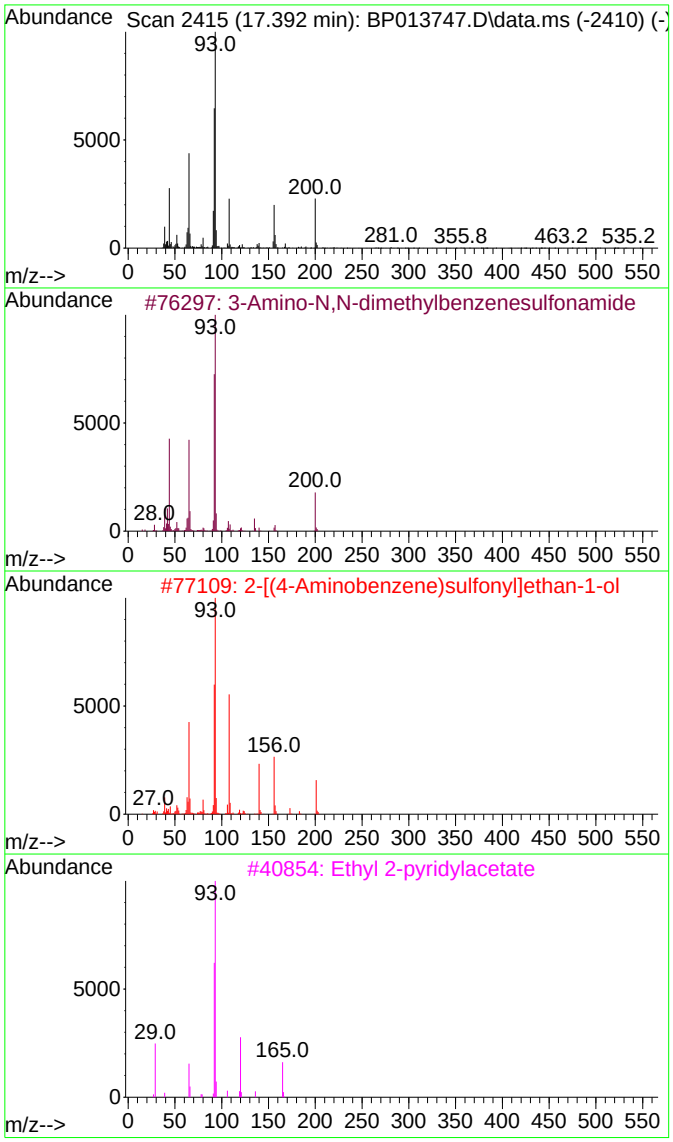
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

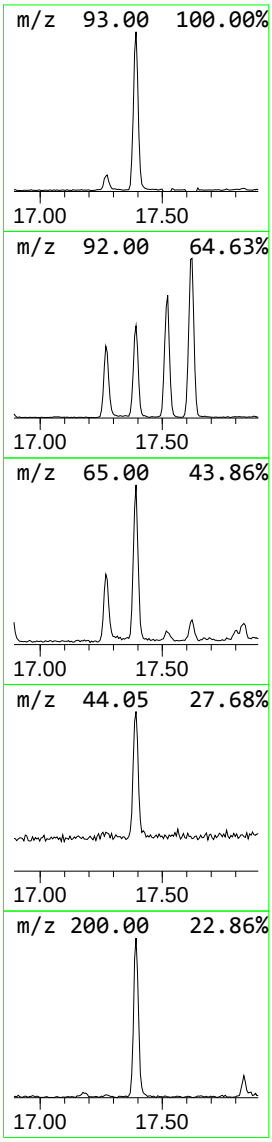
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 3-Amino-N,N-dimethylbenzene... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.392	2.34 ng/ul	438919	Phenanthrene-d10	17.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Amino-N,N-dimethylbenzenesulfo...	200	C8H12N2O2S	006274-18-6	87
2	2-[(4-Aminobenzene)sulfonyl]etha...	201	C8H11NO3S	005246-58-2	72
3	Ethyl 2-pyridylacetate	165	C9H11NO2	002739-98-2	59
4	4-Aminophenyl chlorodifluorometh...	241	C7H6ClF2NO2S	074171-46-3	47
5	3-Picoline, 2-nitro-	138	C6H6N2O2	018368-73-5	46





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Acq On : 03 Feb 2023 18:57
Operator : CG/JU
Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

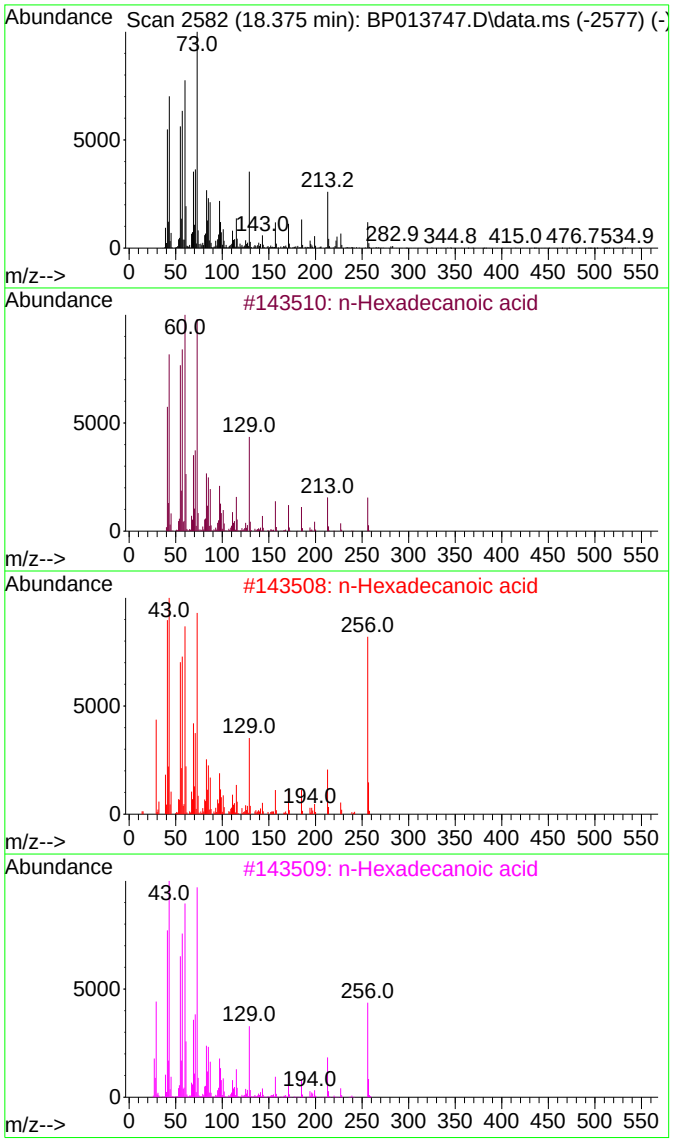
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

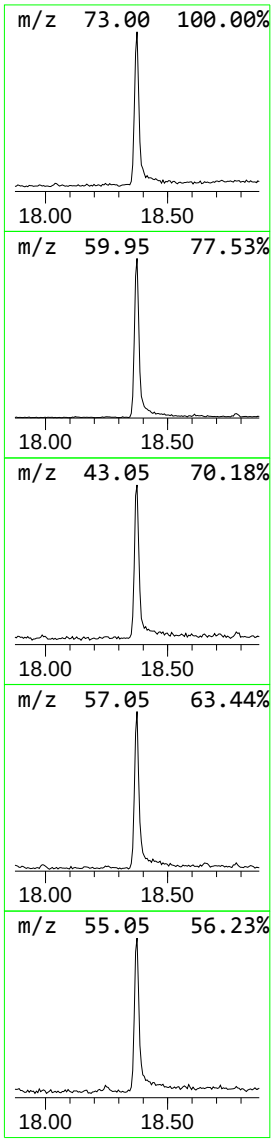
****

Peak Number 15 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.375	3.28 ng/ul	614606	Phenanthrene-d10	17.522

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	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96





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Acq On : 03 Feb 2023 18:57
Operator : CG/JU
Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

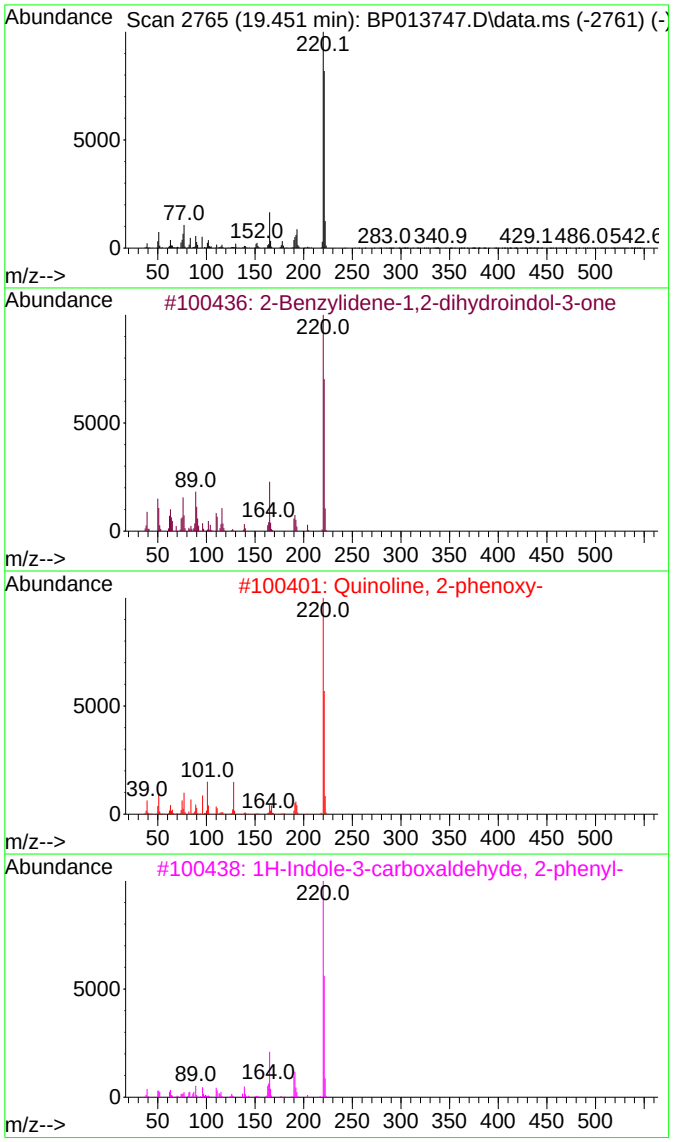
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

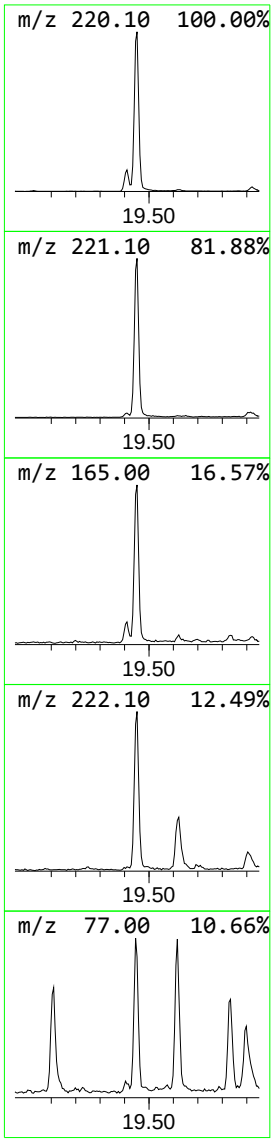
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 2-Benzylidene-1,2-dihydroin... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.451	5.44 ng/ul	1020180	Phenanthrene-d10	17.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Benzylidene-1,2-dihydroindol-3...	221	C15H11NO	039830-75-6	91
2	Quinoline, 2-phenoxy-	221	C15H11NO	040515-82-0	87
3	1H-Indole-3-carboxaldehyde, 2-ph...	221	C15H11NO	025365-71-3	78
4	1H-Quinolin-2-one, 3-phenyl-	221	C15H11NO	038035-81-3	76
5	4,6-Difluoro-2-(3-methylphenylam...	221	C11H9F2N3	1000070-52-7	64





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Acq On : 03 Feb 2023 18:57
Operator : CG/JU
Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

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

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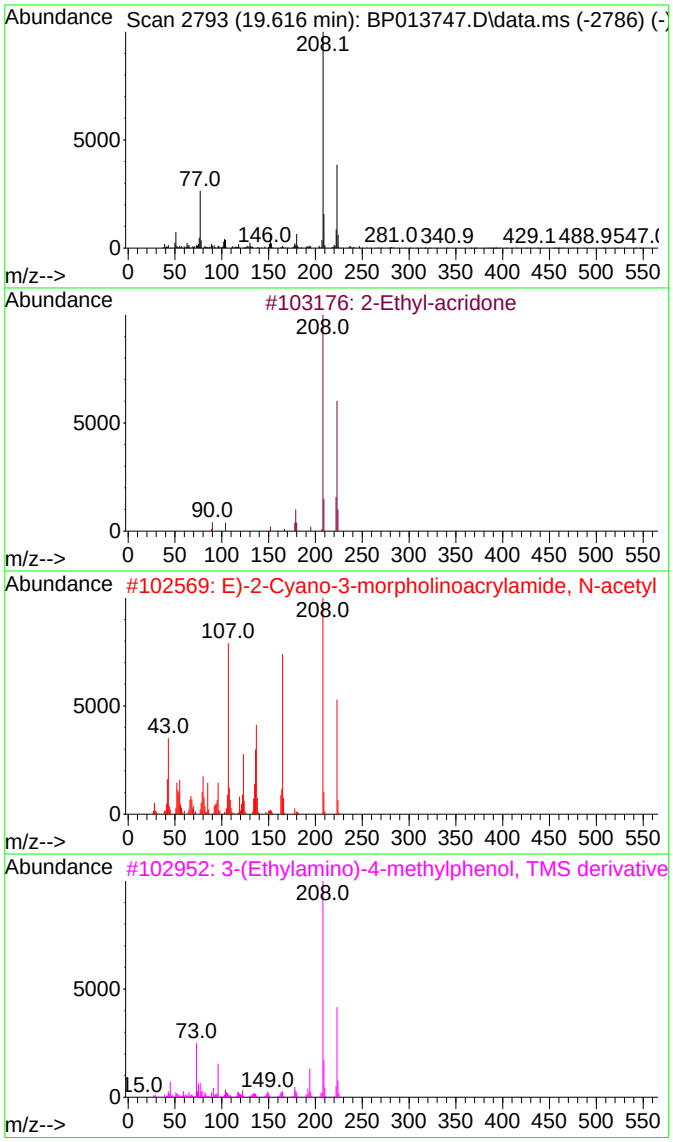
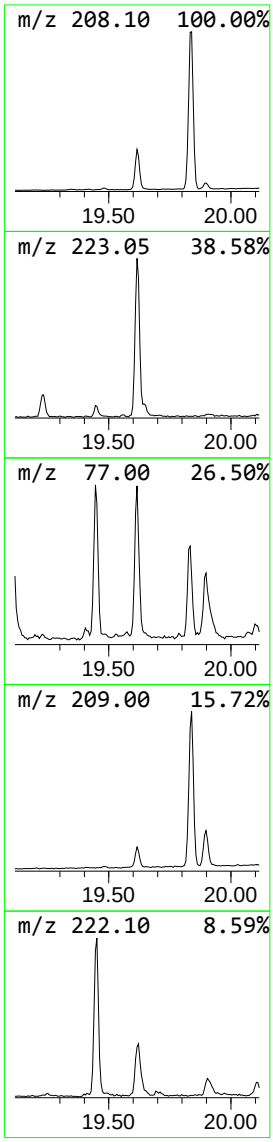
Peak Number 17 2-Ethyl-acridone Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.616	2.50 ng/ul	621410	Chrysene-d12	21.598

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Ethyl-acridone	223	C15H13NO	065753-77-7	78
2	E)-2-Cyano-3-morpholinoacrylamid...	223	C10H13N3O3	1000514-05-3	64
3	3-(Ethylamino)-4-methylphenol, T...	223	C12H21NOSi	1000488-54-1	64
4	Indole, 5-methyl-2-(4-pyridyl)-	208	C14H12N2	107919-92-6	64
5	4-Quinolinol,4-ethenyl-1-ethylde...	223	C14H25NO	038463-58-0	64





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Operator : CG/JU
Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

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

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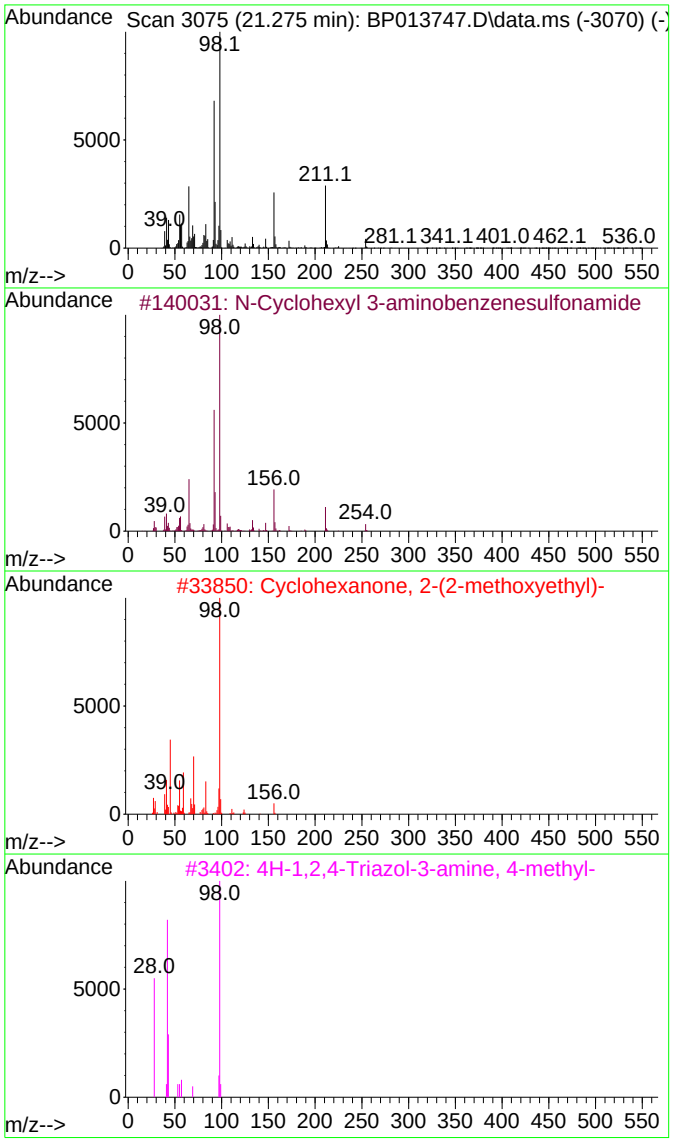
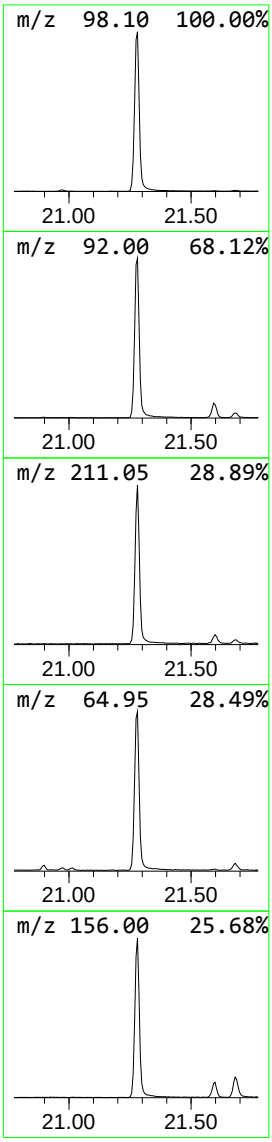
Peak Number 18 N-Cyclohexyl 3-aminobenzene... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.275	20.03 ng/ul	4980420	Chrysene-d12	21.598

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-Cyclohexyl 3-aminobenzenesulfo...	254	C12H18N2O2S	061886-26-8	98
2	Cyclohexanone, 2-(2-methoxyethyl)-	156	C9H16O2	129786-15-8	38
3	4H-1,2,4-Triazol-3-amine, 4-methyl-	98	C3H6N4	016681-76-8	30
4	2,3-2H-4-Methyl-imidazole-2-one	98	C4H6N2O	039799-77-4	30
5	1,3-Cyclohexanediol, trans-	116	C6H12O2	005515-64-0	30





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013747.D
Acq On : 03 Feb 2023 18:57
Operator : CG/JU
Sample : 01381-13
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.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
m-Chloroaniline	9.975	3.0	ng/ul	256414	2	10.957	1732200	20.0
1,3-Pentanediol...	10.240	3.2	ng/ul	278450	2	10.957	1732200	20.0
Benzene, 1,3,5-...	10.822	52.8	ng/ul	4571600	2	10.957	1732200	20.0
Aniline, N-methyl-	12.469	5.8	ng/ul	500593	2	10.957	1732200	20.0
Benzenamine, 2,...	12.975	15.6	ng/ul	2084330	3	14.769	2671670	20.0
Benzophenone	16.122	21.3	ng/ul	2839670	3	14.769	2671670	20.0
Benzenesulfonam...	16.345	6.0	ng/ul	1131490	4	17.522	3749930	20.0
3-Amino-N,N-dim...	17.392	2.3	ng/ul	438919	4	17.522	3749930	20.0
n-Hexadecanoic ...	18.375	3.3	ng/ul	614606	4	17.522	3749930	20.0
2-Benzylidene-1...	19.451	5.4	ng/ul	1020180	4	17.522	3749930	20.0
2-Ethyl-acridone	19.616	2.5	ng/ul	621410	5	21.598	4973500	20.0
N-Cyclohexyl 3-...	21.275	20.0	ng/ul	4980420	5	21.598	4973500	20.0