LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BHA38

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-BP101122.M
Title : SVOA CALIBRATION

Signal : TIC: BP012122.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.070	11	14	32	rVB	596734	853287	9.29%	0.738%
2	3.617	103	107	117	rVB	204943	272409	2.97%	0.236%
3	3.693	117	120	133	rBV	946044	1465130	15.96%	1.267%
4	3.840	142	145	154	rVB	121621	154976	1.69%	0.134%
5	4.017	169	175	192	rVB	6329651	8837682	96.26%	7.645%
6	5.152	356	368	383	rBV	1184905	1834643	19.98%	1.587%
7	5.475	417	423	433	rBV	2619769	5310467	57.84%	4.594%
8	5.558	433	437	443	rVB2	77384	116121	1.26%	0.100%
9	5.764	459	472	480	rBV3	40557	109540	1.19%	0.095%
10	5.846	480	486	499	rVB	1666942	2607274	28.40%	2.255%
11	6.511	589	599	606	rBV3	116196	216539	2.36%	0.187%
12	6.822	641	652	664	rBV5	45573	119020	1.30%	0.103%
13	6.928	664	670	674	rBV	225197	346367	3.77%	0.300%
14	6.993	674	681	682	rVV3	119111	174989	1.91%	0.151%
15	7.017	682	685	689	rVV	259487	440901	4.80%	0.381%
16	7.064	689	693	697	rVB2	234173	393896	4.29%	0.341%
17	7.111	697	701	706	rVB3	75580	113499	1.24%	0.098%
18	7.181	706	713	718	rBV	1053764	1647231	17.94%	1.425%
19	7.228	718	721	725	rVB	220916	297070	3.24%	0.257%
20	7.275	725	729	737	rVB	192175	294772	3.21%	0.255%
21	7.381	737	747	758	rBV	950163	1733988	18.89%	1.500%
22	7.487	758	765	771	rVB	423475	705565	7.68%	0.610%
23	7.552	771	776	781	rVB3	106973	165366	1.80%	0.143%
24	7.846	816	826	831	rBV	828330	1412712	15.39%	1.222%
25	7.958	839	845	855	rBV	353230	576000	6.27%	0.498%
26	8.046	855	860	865	rBV	164370	231638	2.52%	0.200%
27	8.216	881	889	895	rBV2	524436	1037413	11.30%	0.897%
28	8.287	895	901	908	rBV2	1080163	2265893	24.68%	1.960%
29	8.393	915	919	926	rVB	78138	116549	1.27%	0.101%
30	8.475	926	933	941	rBV	1964962	3174895	34.58%	2.746%
31	8.563	941	948	954	rVB	751618	1174944	12.80%	1.016%
32	8.634	954	960	962	rBV2	124056	228087	2.48%	0.197%
33	8.852	991	997	1004	rVB3	77655	147883	1.61%	0.128%
34	8.952	1004	1014	1019	rBV	157056	249803	2.72%	0.216%
35	9.016	1019	1025	1034	rBV	1028861	1912645	20.83%	1.654%
36	9.081	1034	1036	1046	rVV4	121296	226115	2.46%	0.196%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BHA38

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-BP101122.M
Title : SVOA CALIBRATION

37	9.169	1046	1051	1061	rVB	343437	530738	5.78%	0.459%
38	9.281	1062	1070	1079	rVB5	48677	116144	1.26%	0.100%
39	9.475	1097	1103	1109	rVB	93593	156677	1.71%	0.136%
40	9.534	1109	1113	1120	rBV	72486	114925	1.25%	0.099%
41	9.640	1124	1131	1136	rBV2	87418	166359	1.81%	0.144%
42	9.740	1136	1148	1160	rBV	832143	1622417	17.67%	1.403%
43	9.899	1171	1175	1180	rVB	93080	133129	1.45%	0.115%
44	10.069	1197	1204	1212	rVV5	195415	508052	5.53%	0.439%
45	10.146	1212	1217	1222	rVV3	71439	113185	1.23%	0.098%
46	10.275	1229	1239	1248	rBV	1246812	2294913	25.00%	1.985%
47	10.469	1268	1272	1277	rVB2	74045	110931	1.21%	0.096%
48	10.657	1296	1304	1308	rVV	1094556	1860067	20.26%	1.609%
49	10.705	1308	1312	1316	rVV	376878	656638	7.15%	0.568%
50	10.746	1316	1319	1322	rVV	201637	333062	3.63%	0.288%
51	10.805	1322	1329	1343	rVB2	4907459	9181362	100.00%	7.942%
52	11.069	1359	1374	1379	rBV5	129545	382578	4.17%	0.331%
53	11.257	1399	1406	1414	rVV4	121918	304438	3.32%	0.263%
54	11.675	1473	1477	1483	rVV4	59555	117476	1.28%	0.102%
55	11.740	1483	1488	1494	rVV	261538	412859	4.50%	0.357%
56	11.999	1522	1532	1538	rBV	530912	1372365	14.95%	1.187%
57	12.316	1578	1586	1597	rVB2	179161	329139	3.58%	0.285%
58	12.475	1599	1613	1615	rVV8	59121	167888	1.83%	0.145%
59	12.534	1619	1623	1631	rVB	260144	427916	4.66%	0.370%
60	13.093	1713	1718	1721	rVV3	84136	139438	1.52%	0.121%
61	13.134	1721	1725	1731	rVB	149633	210245	2.29%	0.182%
62	13.910	1851	1857	1864	rVV	2063388	2861021	31.16%	2.475%
63	13.975	1864	1868	1877	rVB	261396	420933	4.58%	0.364%
64	14.193	1899	1905	1914	rVV	2530643	3801405	41.40%	3.288%
65	14.287	1917	1921	1935	rVB3	48662	124358	1.35%	0.108%
66	14.498	1949	1957	1965	rBV2	1616542	2805865	30.56%	2.427%
67	14.557	1965	1967	1976	rVB	138472	204913	2.23%	0.177%
68	14.716	1988	1994	2004	rBV	158678	329815	3.59%	0.285%
69	14.898	2019	2025	2032	rVB	48891	127059	1.38%	0.110%
70	15.493	2120	2126	2133	rBV2	3789226	5968950	65.01%	5.163%
71	15.616	2142	2147	2153	rBV	821688	1102312	12.01%	0.954%
72	16.010	2208	2214	2222	rBV4	234935	451990	4.92%	0.391%
73	16.192	2241	2245	2249	rVV	138063	215663	2.35%	0.187%
74	16.234	2250	2252	2265	rVB10	64875	140768	1.53%	0.122%
75	16.557	2300	2307	2313	rVB	247497	363147	3.96%	0.314%
76	17.004	2378	2383	2389	rVB3	67682	111955	1.22%	0.097%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

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-BP101122.M
Title : SVOA CALIBRATION

77	17.216	2414	2419	2422	rBV	547538	729048	7.94%	0.631%
78	17.257	2422	2426	2437	rVB	1788461	2622052	28.56%	2.268%
79	17.357	2437	2443	2450	rBV	3378289	4994059	54.39%	4.320%
80	17.634	2484	2490	2494	rVB2	72170	116661	1.27%	0.101%
81	17.781	2511	2515	2519	rVV	90468	110934	1.21%	0.096%
82	17.934	2536	2541	2546	rVB3	79215	148367	1.62%	0.128%
83	18.139	2572	2576	2584	rVB	124866	165418	1.80%	0.143%
84	18.286	2597	2601	2609	rBV	271985	353776	3.85%	0.306%
85	18.386	2612	2618	2620	rVV	232643	324711	3.54%	0.281%
86	18.434	2620	2626	2632	rVB	382284	701960	7.65%	0.607%
87	19.110	2738	2741	2747	rVV	163886	227504	2.48%	0.197%
88	19.316	2772	2776	2782	rBV3	162927	206307	2.25%	0.178%
89	19.375	2782	2786	2798	rVV2	123567	211786	2.31%	0.183%
90	19.592	2818	2823	2828	rBV	4748016	6182136	67.33%	5.348%
91	19.639	2828	2831	2841	rVB	692015	867586	9.45%	0.750%
92	19.828	2861	2863	2869	rVB3	109831	141041	1.54%	0.122%
93	20.033	2895	2898	2902	rBV	101062	112076	1.22%	0.097%
94	20.086	2904	2907	2916	rVB7	109622	243095	2.65%	0.210%
95	20.298	2940	2943	2948	rVB	188992	227701	2.48%	0.197%
96	20.451	2966	2969	2973	rBV	129371	158730	1.73%	0.137%
97	21.345	3116	3121	3127	rBV	2664035	3331648	36.29%	2.882%
98	21.469	3138	3142	3146	rBV	523391	689431	7.51%	0.596%
99	23.610	3499	3506	3517	rVB2	3685408	7301536	79.53%	6.316%
100	23.763	3526	3532	3544	rVB2	1837788	3677496	40.05%	3.181%

Sum of corrected areas: 115603463

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012122.D
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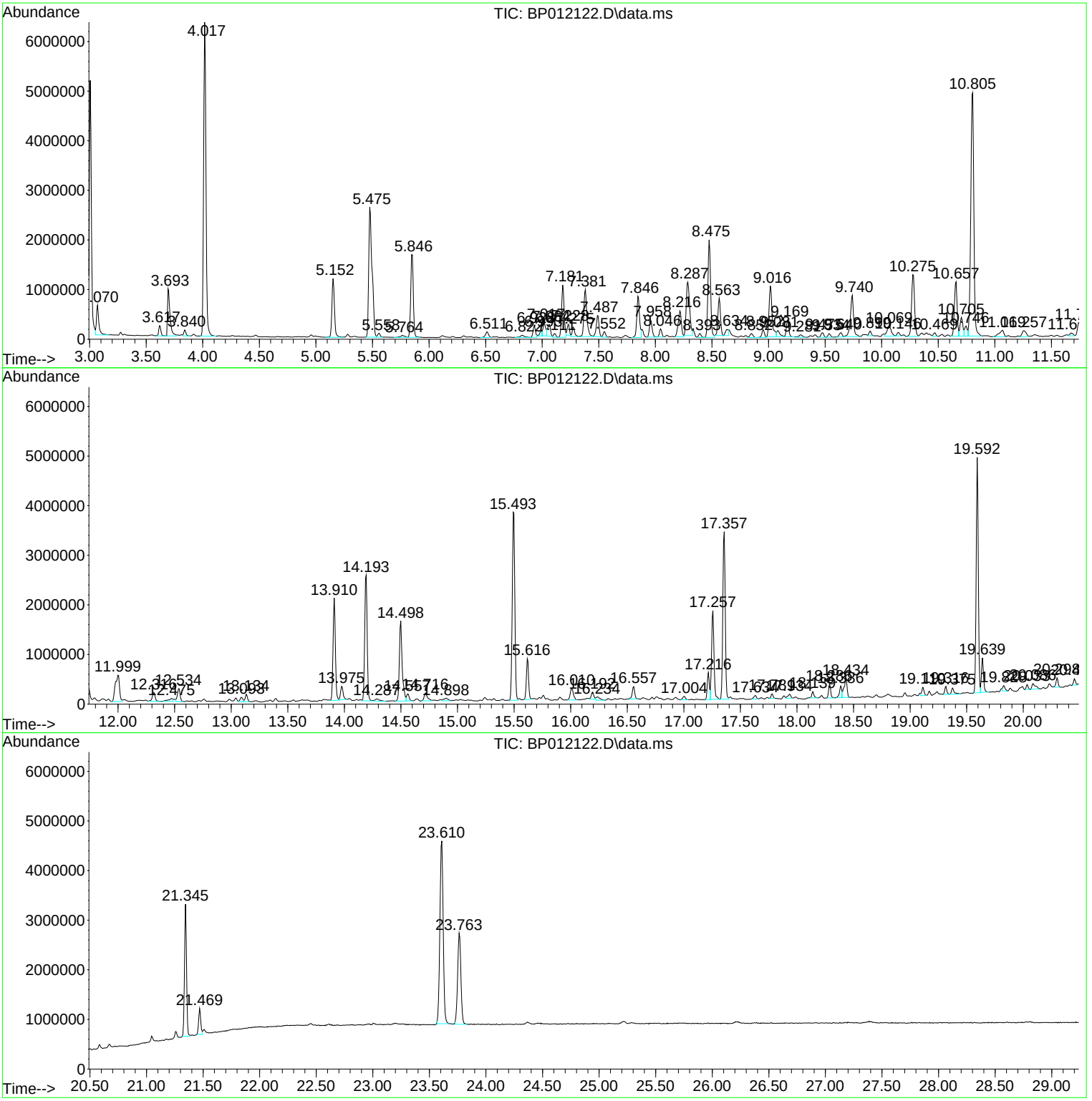
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

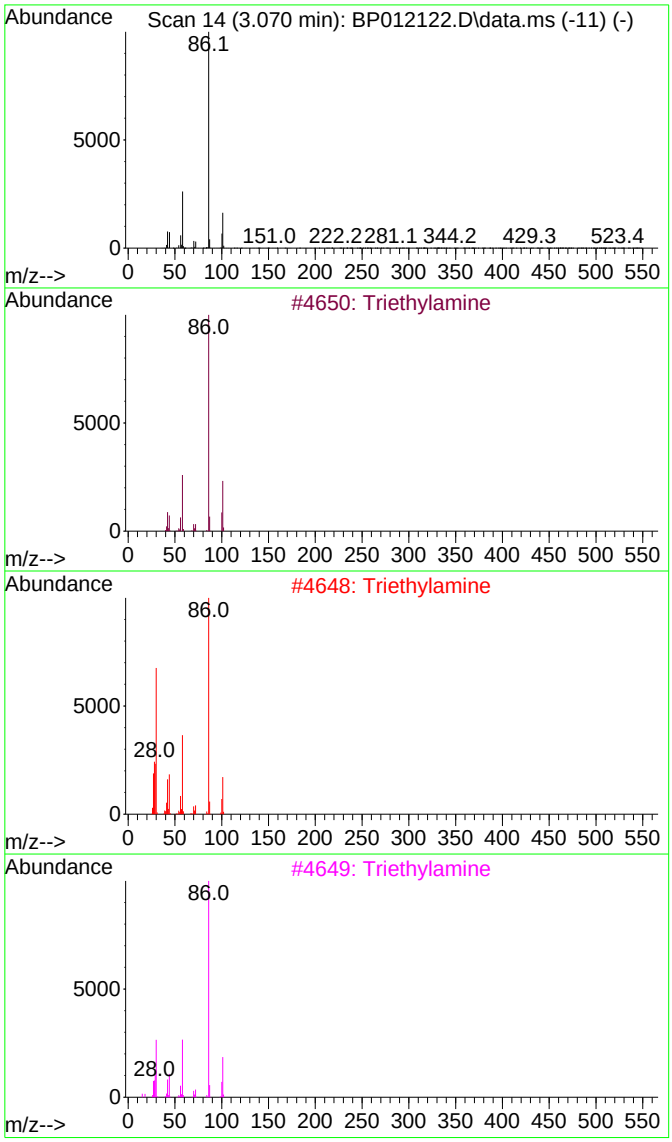
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

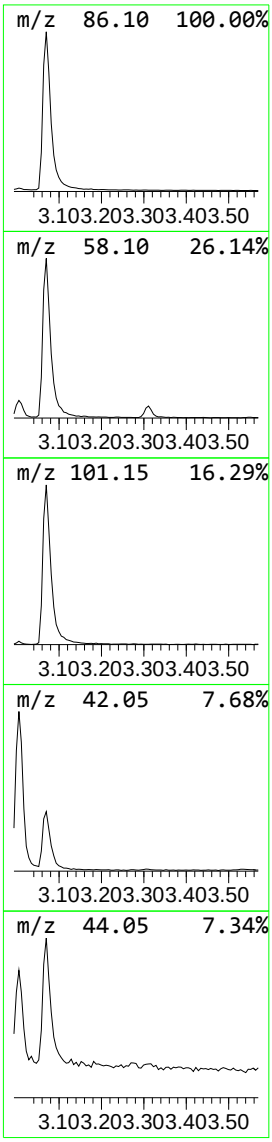
****

Peak Number 1 Triethylamine Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.070	12.08 ng/ul	853287	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triethylamine	101	C6H15N	000121-44-8	91
2	Triethylamine	101	C6H15N	000121-44-8	91
3	Triethylamine	101	C6H15N	000121-44-8	87
4	Triethylamine	101	C6H15N	000121-44-8	87
5	Tetraethylammonium chloride	165	C8H20ClN	000056-34-8	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

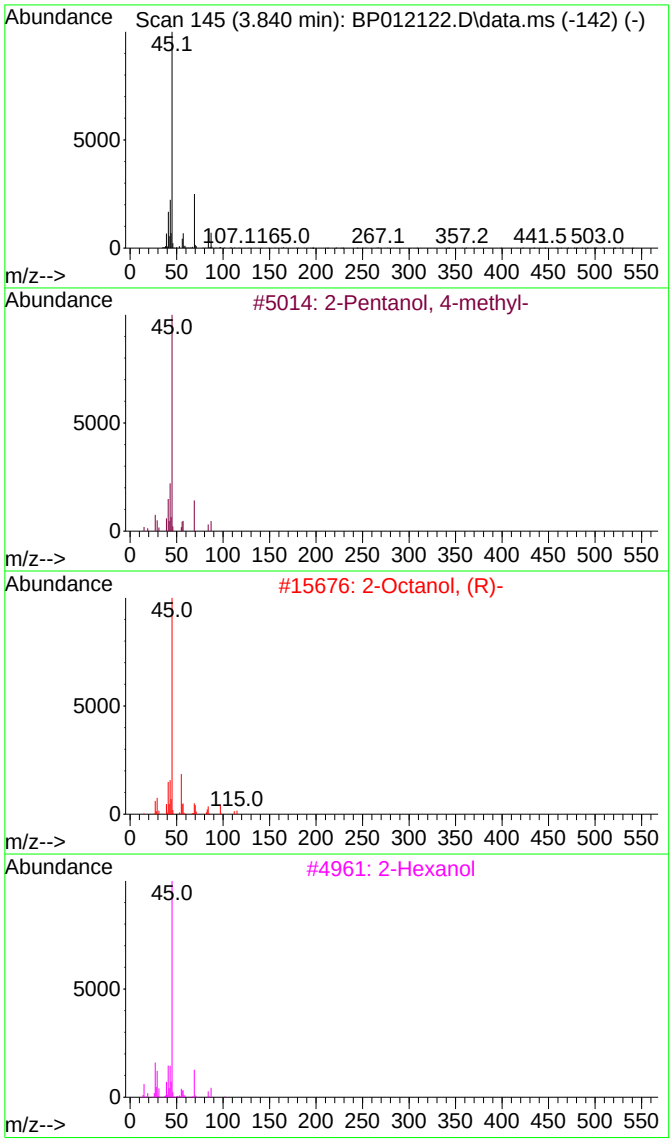
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

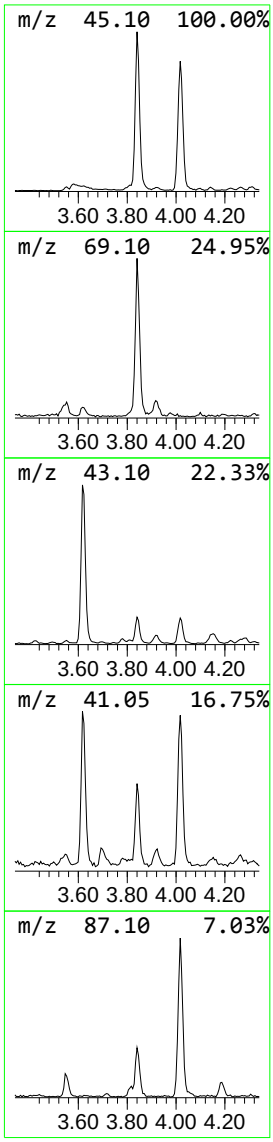
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Pentanol, 4-methyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.840	2.19 ng/ul	154976	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	83
2	2-Octanol, (R)-	130	C8H18O	005978-70-1	78
3	2-Hexanol	102	C6H14O	000626-93-7	64
4	(+-)-5-Methyl-2-hexanol	116	C7H16O	111768-09-3	64
5	2-Hexanol	102	C6H14O	000626-93-7	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

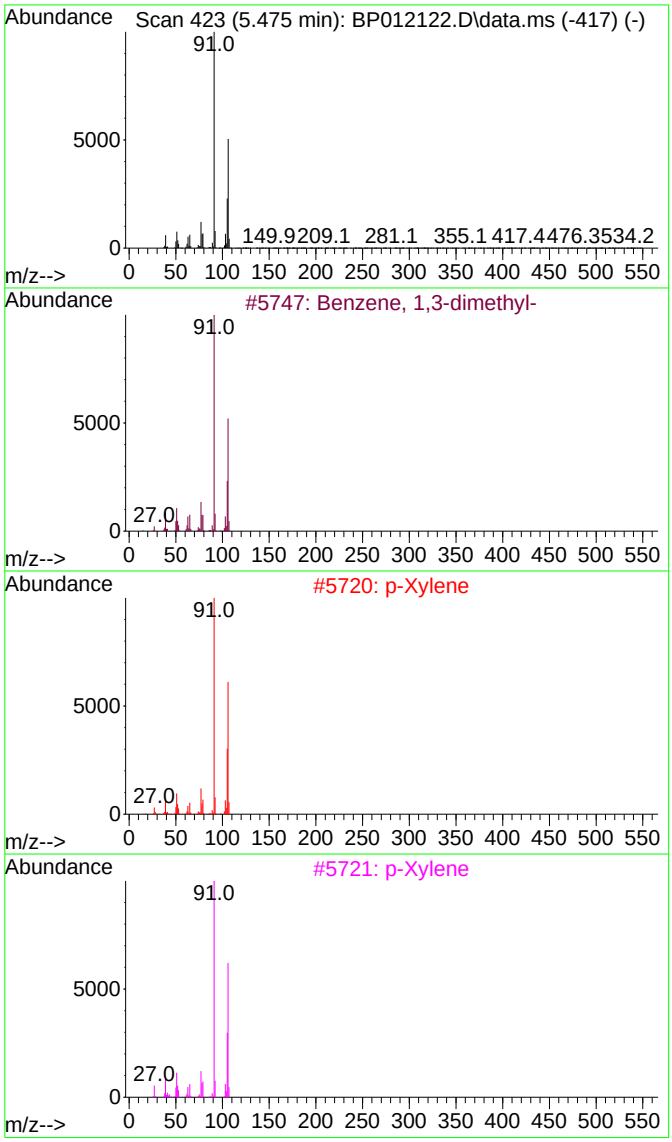
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

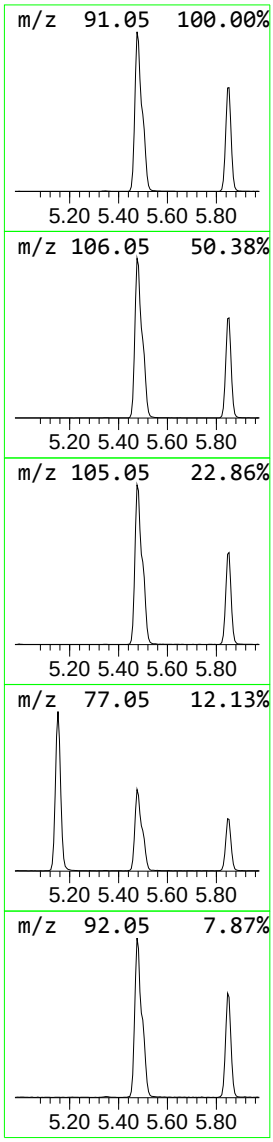
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.475	75.18 ng/ul	5310470	1,4-Dichlorobenzene-d4	7.846

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





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

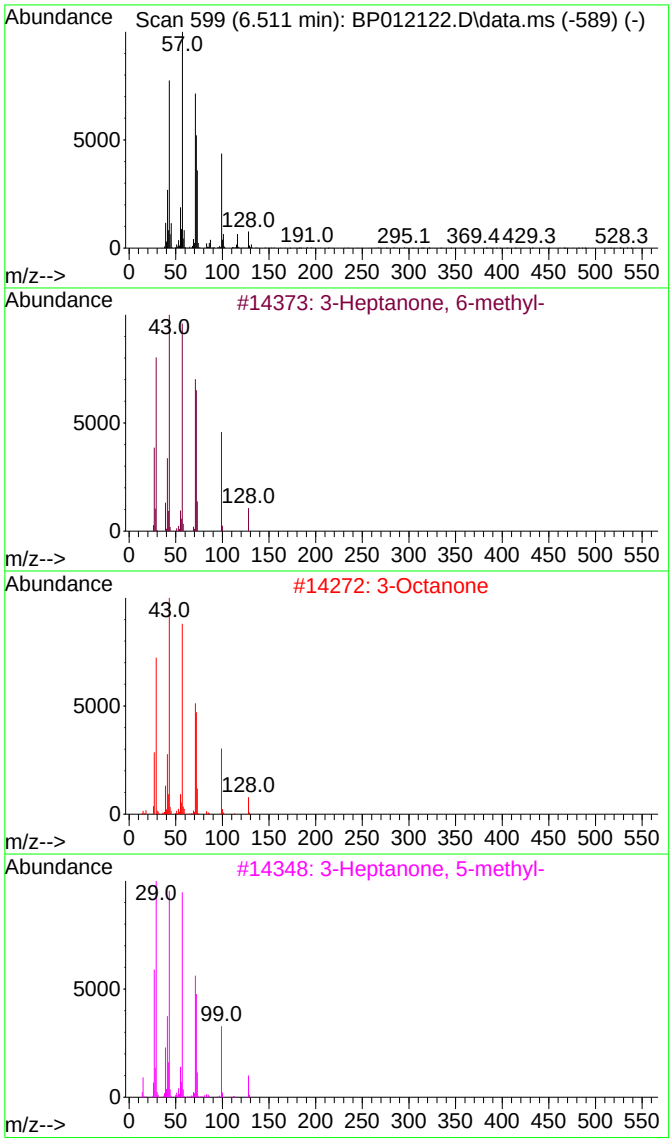
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

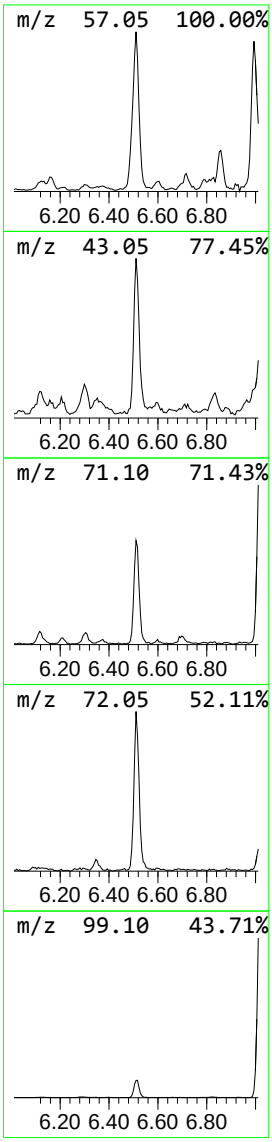
****

Peak Number 8 3-Heptanone, 6-methyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.511	3.07 ng/ul	216539	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptanone, 6-methyl-	128	C8H16O	000624-42-0	87
2	3-Octanone	128	C8H16O	000106-68-3	80
3	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	72
4	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	64
5	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

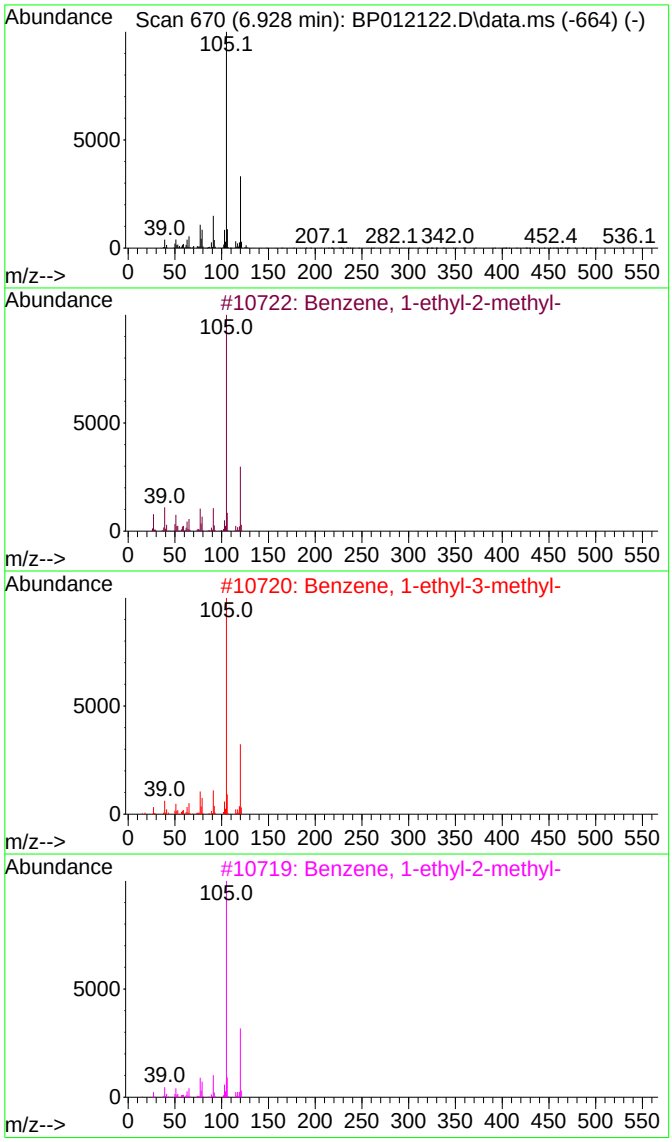
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

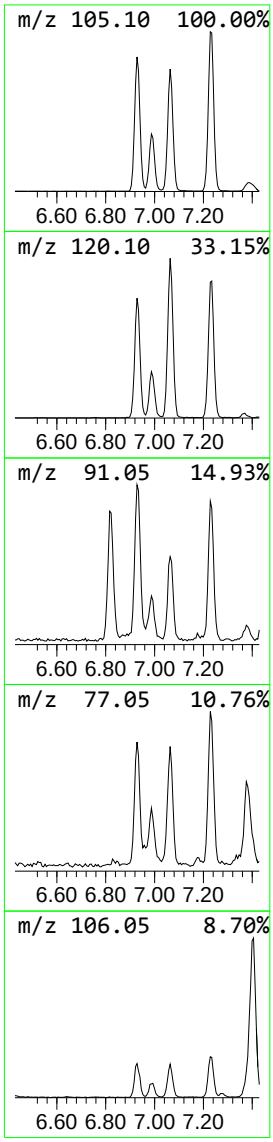
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.928	4.90 ng/ul	346367	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90





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Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

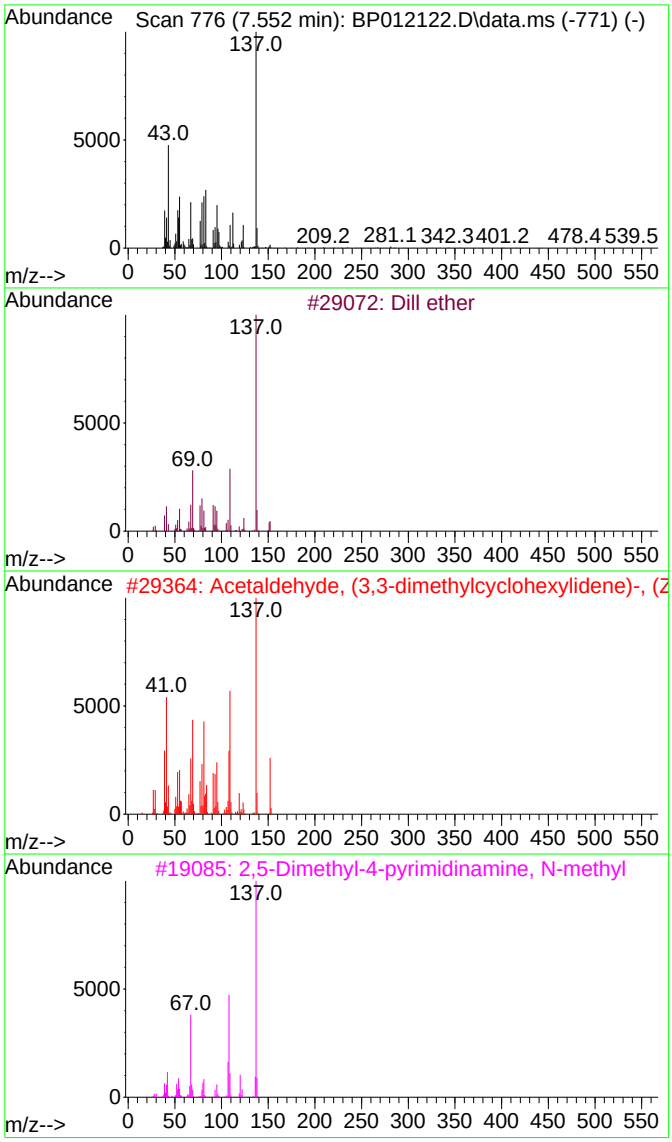
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

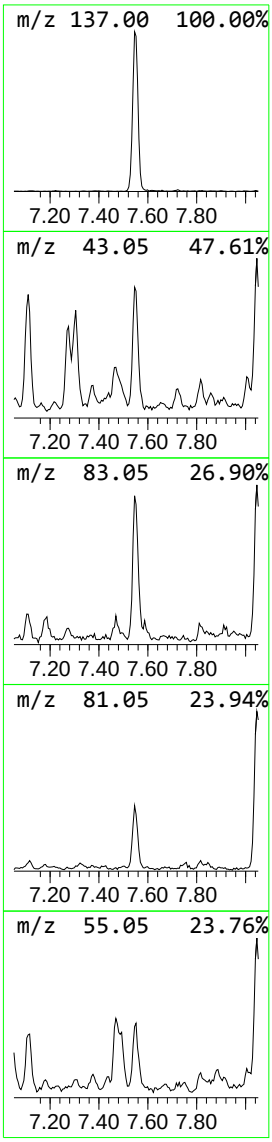
****

Peak Number 11 Dill ether Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.552	2.34 ng/ul	165366	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dill ether	152	C10H16O	074410-10-9	53
2	Acetaldehyde, (3,3-dimethylcyclo...	152	C10H16O	026532-24-1	45
3	2,5-Dimethyl-4-pyrimidinamine, N...	137	C7H11N3	052698-60-9	43
4	Naphthalene, 1,1'-ethylidenebis[...	302	C22H38	054934-70-2	40
5	3,5-Diisopropylpyrazole	152	C9H16N2	017536-00-4	38





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Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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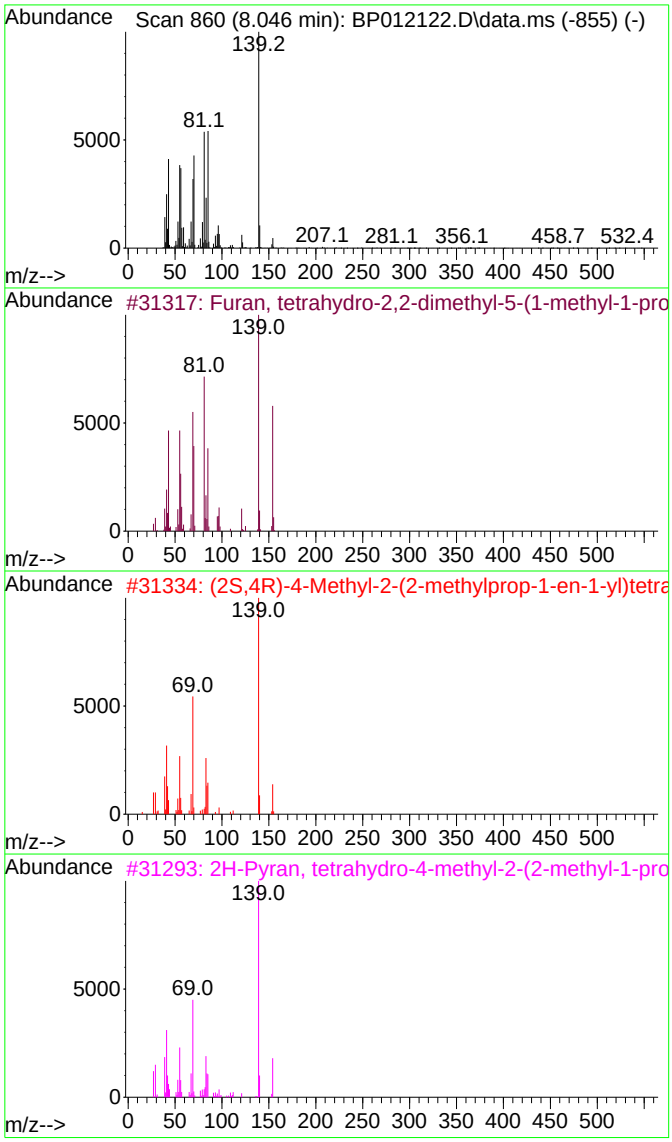
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

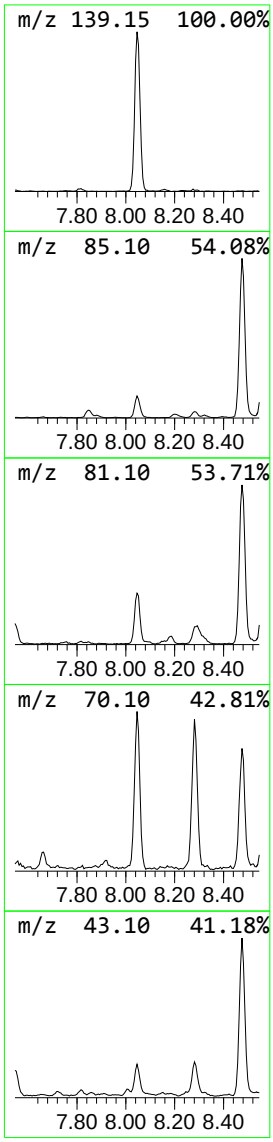
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Furan, tetrahydro-2,2-dimet... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.046	3.28 ng/ul	231638	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Furan, tetrahydro-2,2-dimethyl-5...	154	C10H18O	007416-35-5	50
2	(2S,4R)-4-Methyl-2-(2-methylprop...	154	C10H18O	003033-23-6	38
3	2H-Pyran, tetrahydro-4-methyl-2-...	154	C10H18O	016409-43-1	32
4	(2S,4R)-4-Methyl-2-(2-methylprop...	154	C10H18O	003033-23-6	32
5	Cyclobutylamine, N-trifluoroacetyl-	167	C6H8F3NO	1000461-81-7	32





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Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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

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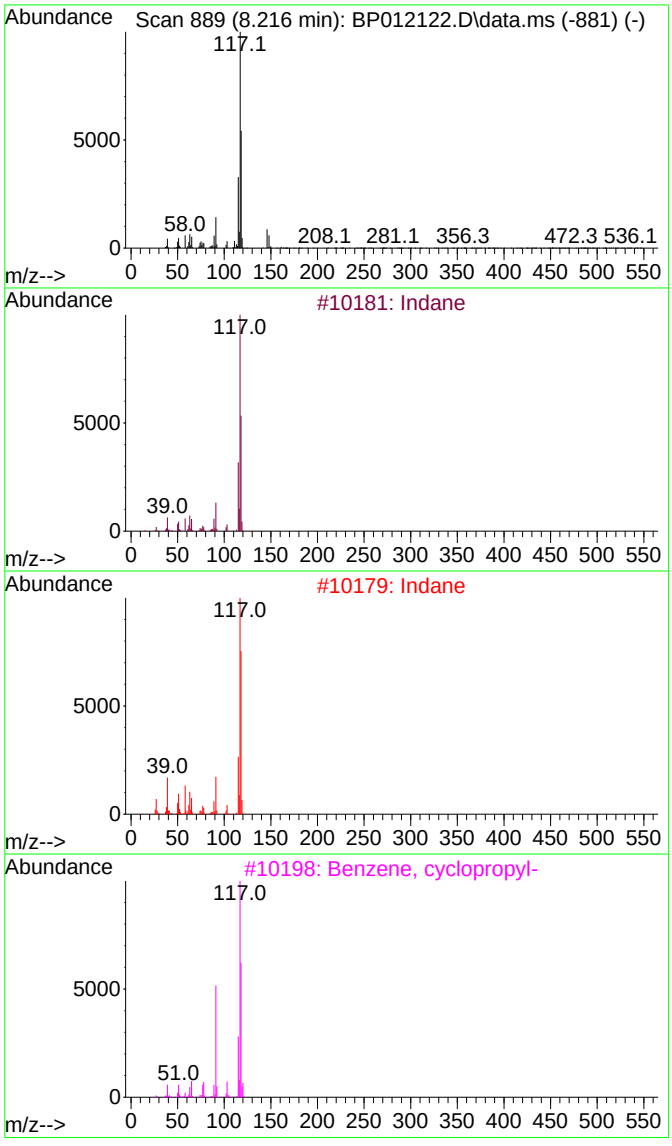
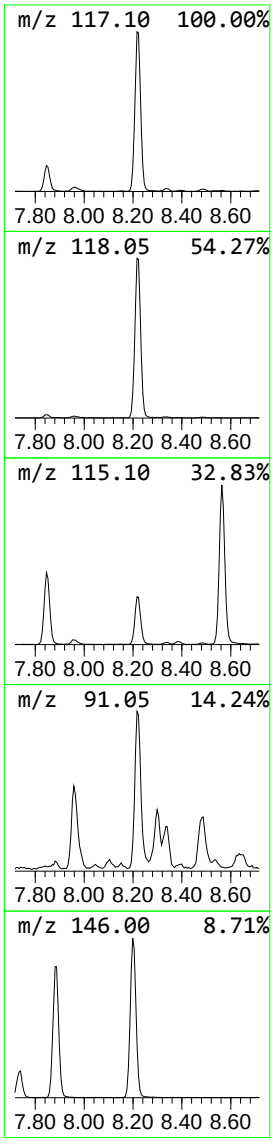
Peak Number 14 Indane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.216	14.69 ng/ul	1037410	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	90
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
4	Deltacyclene	118	C9H10	007785-10-6	72
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72





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Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

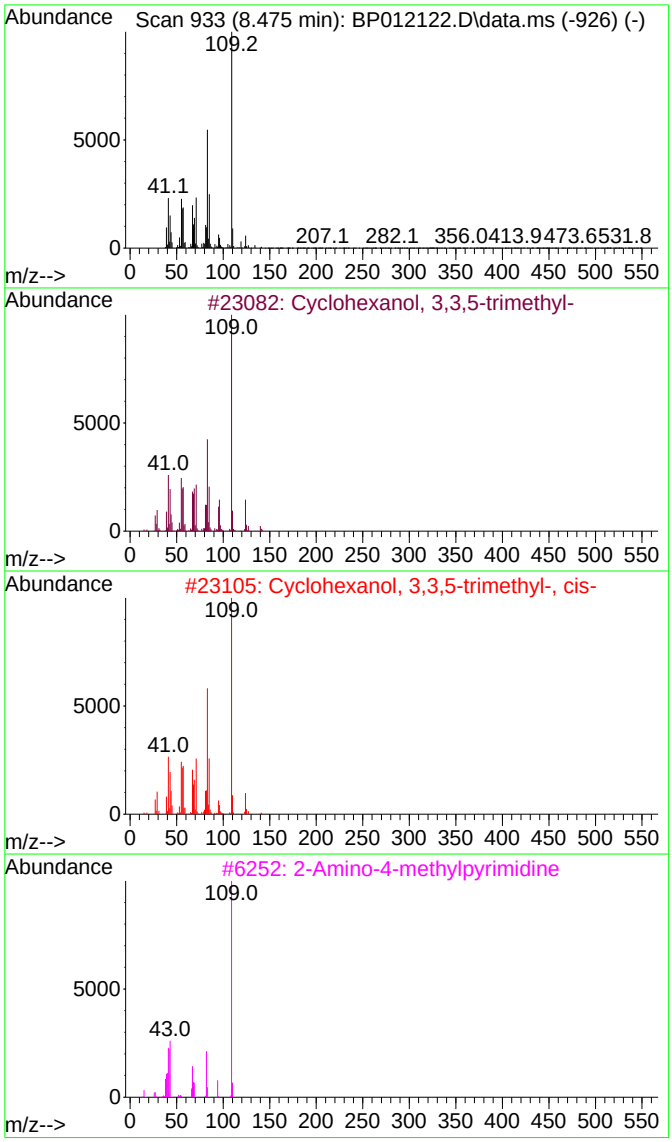
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

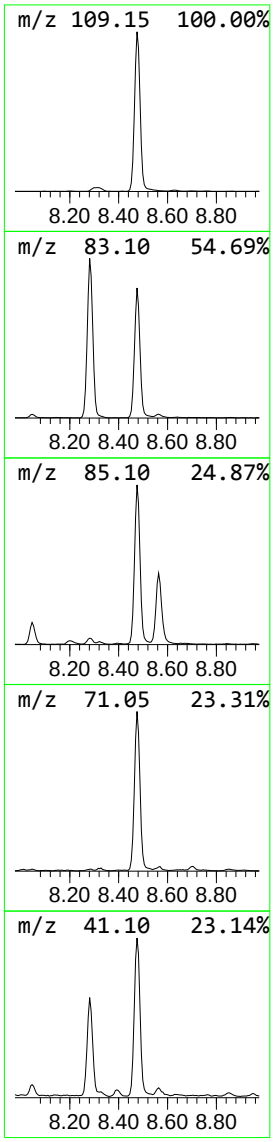
****

Peak Number 15 Cyclohexanol, 3,3,5-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.475	44.95 ng/ul	3174900	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	86
2	Cyclohexanol, 3,3,5-trimethyl-, ...	142	C9H18O	000933-48-2	43
3	2-Amino-4-methylpyrimidine	109	C5H7N3	000108-52-1	43
4	Cyclohexanol, 3,3,5-trimethyl-, ...	184	C11H20O2	024691-16-5	43
5	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	38





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Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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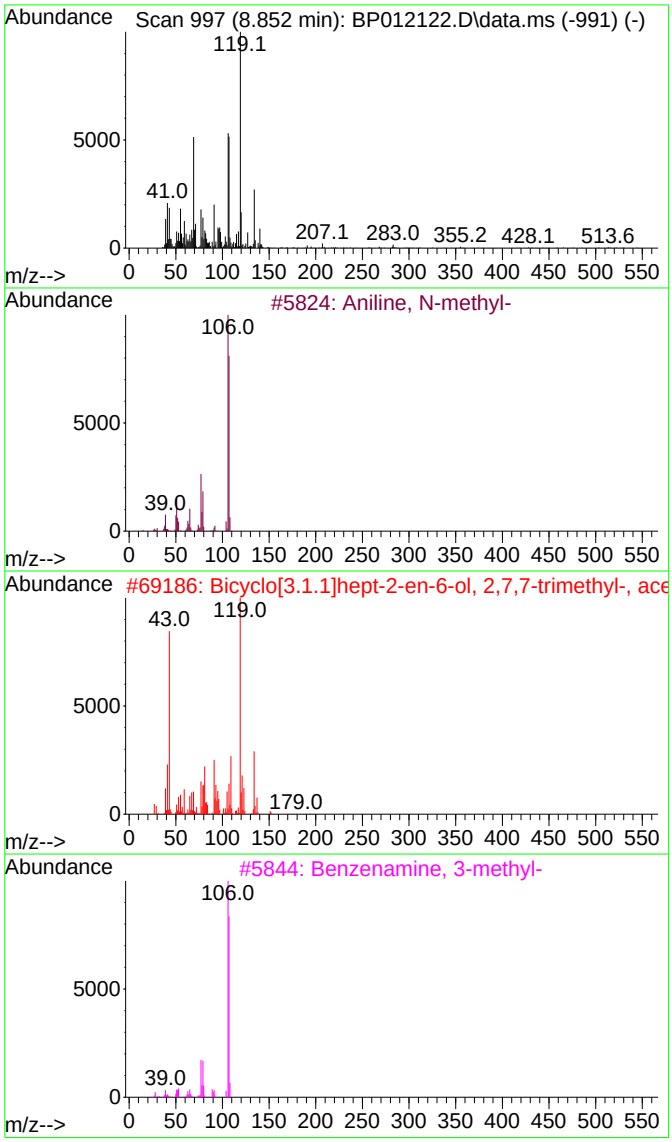
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

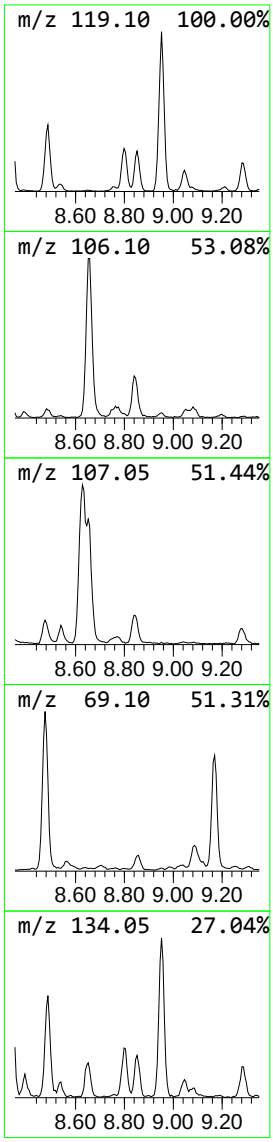
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Aniline, N-methyl- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.852	2.09 ng/ul	147883	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Aniline, N-methyl-	107	C7H9N	000100-61-8	74
2	Bicyclo[3.1.1]hept-2-en-6-ol, 2,...	194	C12H18O2	050764-55-1	49
3	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	47
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	46
5	o-Cymene	134	C10H14	000527-84-4	46





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Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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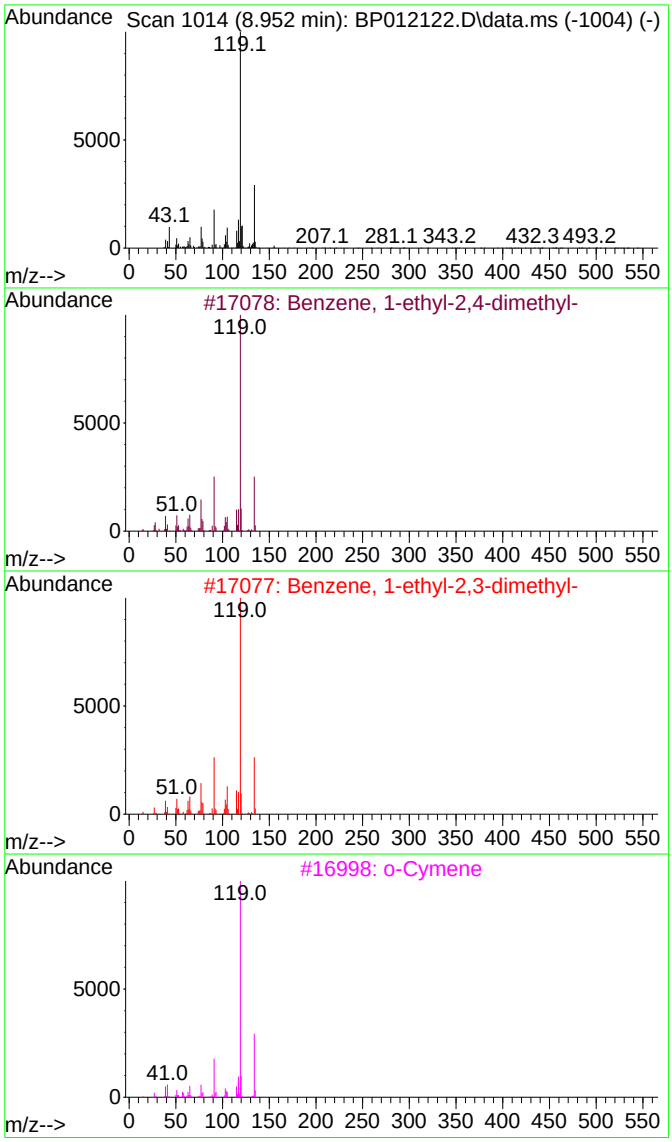
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

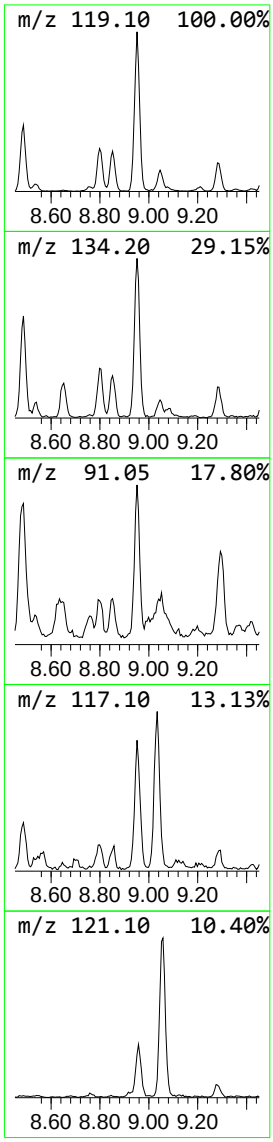
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.952	3.54 ng/ul	249803	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
3	o-Cymene	134	C10H14	000527-84-4	93
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93
5	p-Cymene	134	C10H14	000099-87-6	93





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Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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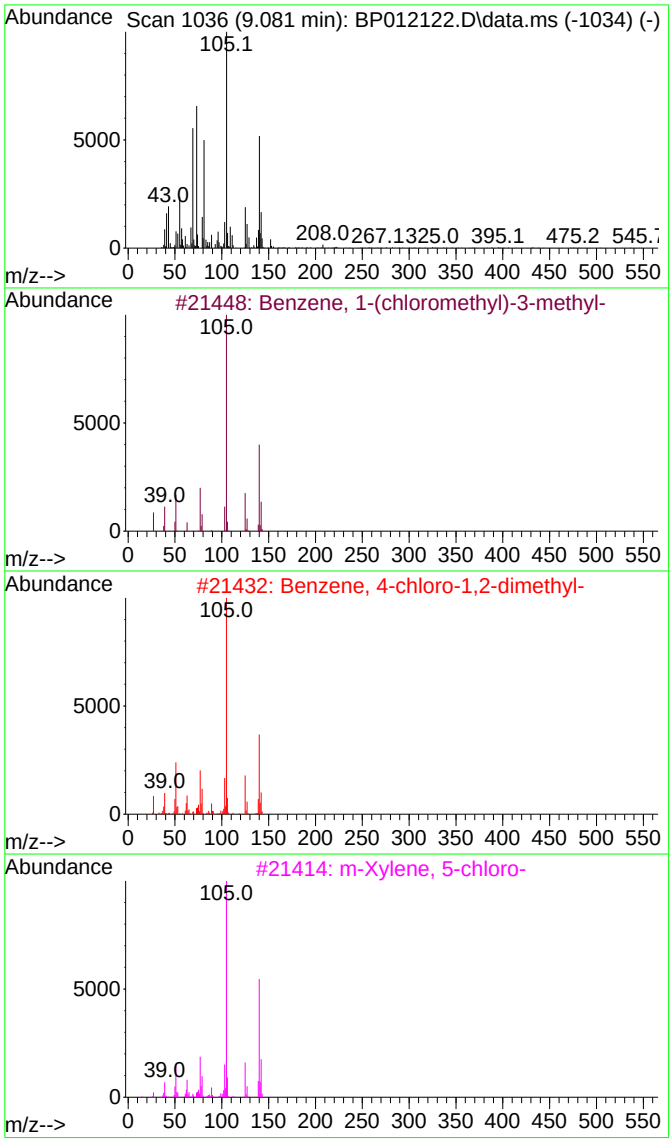
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

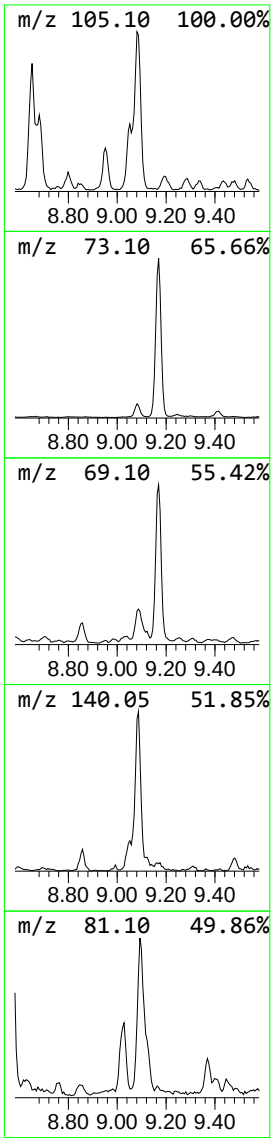
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzene, 1-(chloromethyl)-3... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.081	3.20 ng/ul	226115	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-(chloromethyl)-3-methyl-	140	C8H9Cl	000620-19-9	68
2	Benzene, 4-chloro-1,2-dimethyl-	140	C8H9Cl	000615-60-1	68
3	m-Xylene, 5-chloro-	140	C8H9Cl	000556-97-8	68
4	Benzene, 4-chloro-1,2-dimethyl-	140	C8H9Cl	000615-60-1	68
5	Benzene, (1-chloroethyl)-	140	C8H9Cl	000672-65-1	59





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Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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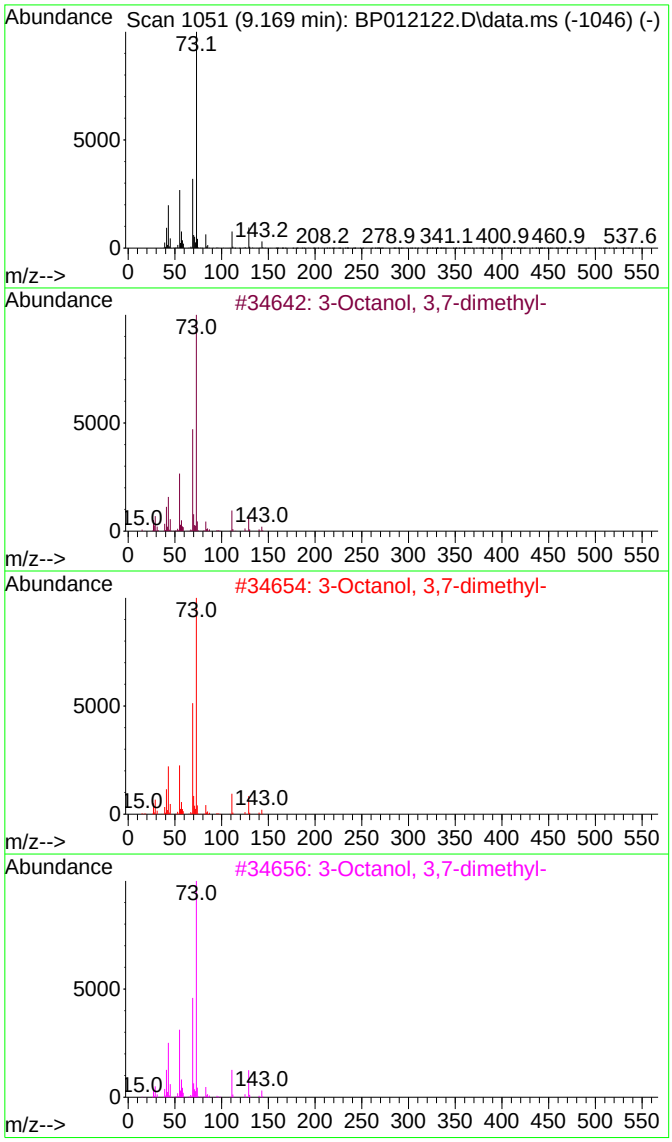
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

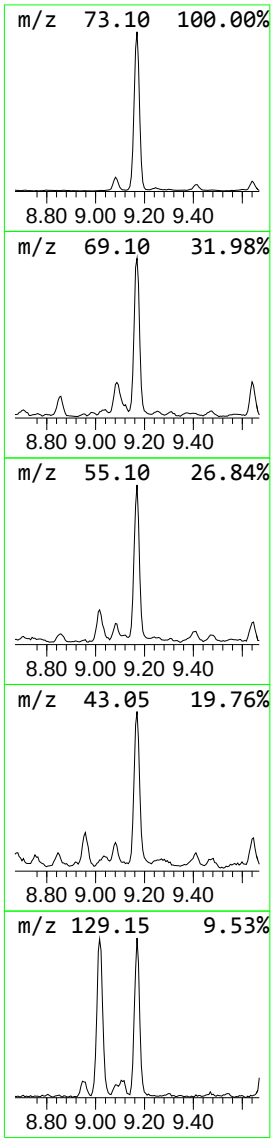
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 3-Octanol, 3,7-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.169	7.51 ng/ul	530738	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	78
2	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	64
3	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	59
4	Decane, 3-methoxy-	172	C11H24O	055955-64-1	53
5	3-Octanol, 3,6-dimethyl-	158	C10H22O	000151-19-9	42





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ALS Vial : 93 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

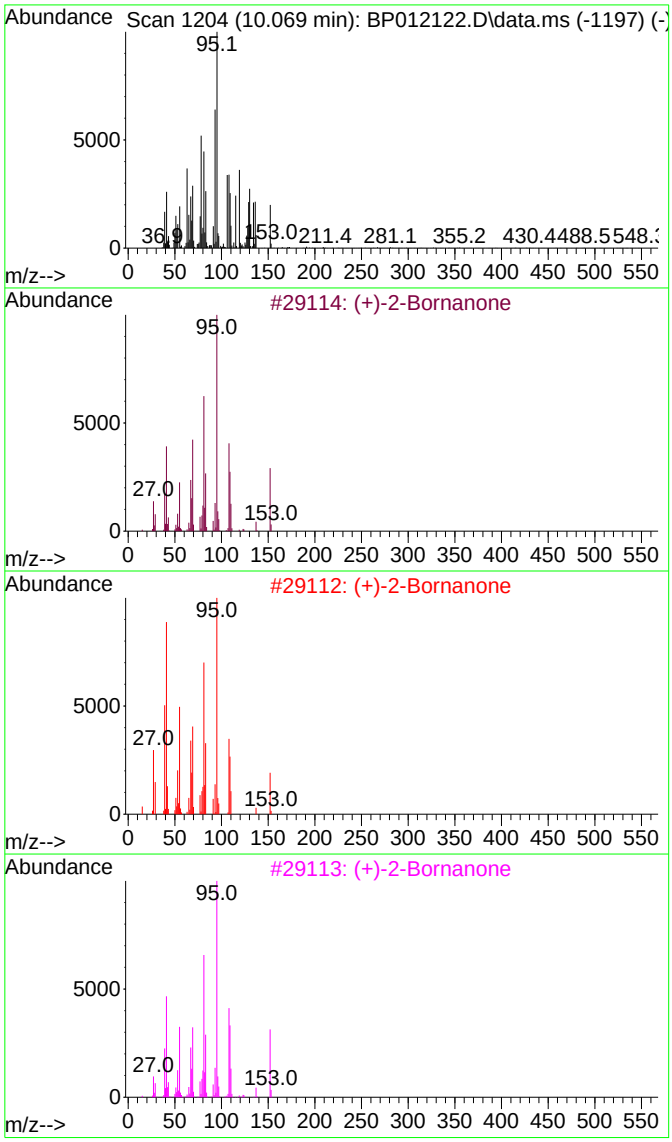
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

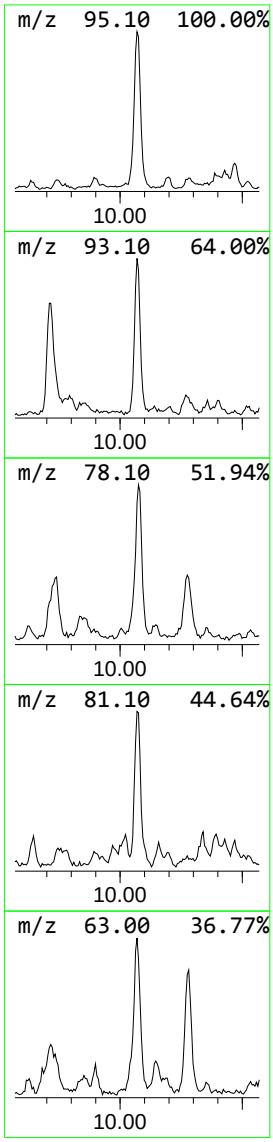
****

Peak Number 20 (+)-2-Bornanone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.069	5.46 ng/ul	508052	Naphthalene-d8	10.658

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-2-Bornanone	152	C10H16O	000464-49-3	95
2	(+)-2-Bornanone	152	C10H16O	000464-49-3	87
3	(+)-2-Bornanone	152	C10H16O	000464-49-3	74
4	(+)-2-Bornanone	152	C10H16O	000464-49-3	70
5	(+)-2-Bornanone	152	C10H16O	000464-49-3	70





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Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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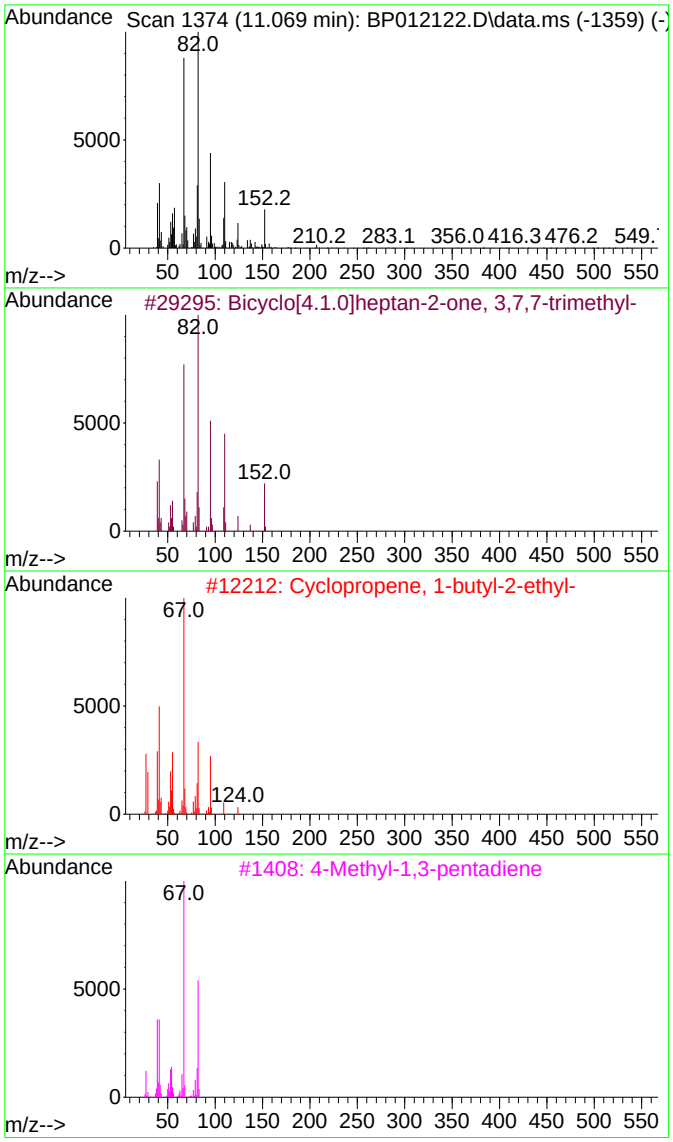
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

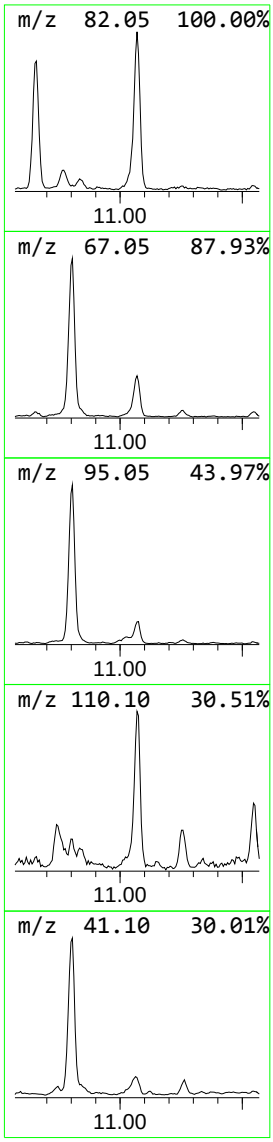
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Bicyclo[4.1.0]heptan-2-one,... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.069	4.11 ng/ul	382578	Naphthalene-d8	10.658

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.1.0]heptan-2-one, 3,7,...	152	C10H16O	000497-62-1	94
2	Cyclopropene, 1-butyl-2-ethyl-	124	C9H16	050915-91-8	58
3	4-Methyl-1,3-pentadiene	82	C6H10	000926-56-7	49
4	4-Nonyne	124	C9H16	020184-91-2	49
5	4-Hexen-1-ol, acetate	142	C8H14O2	072237-36-6	47





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ALS Vial : 93 Sample Multiplier: 1

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Quant Title : SVOA CALIBRATION

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

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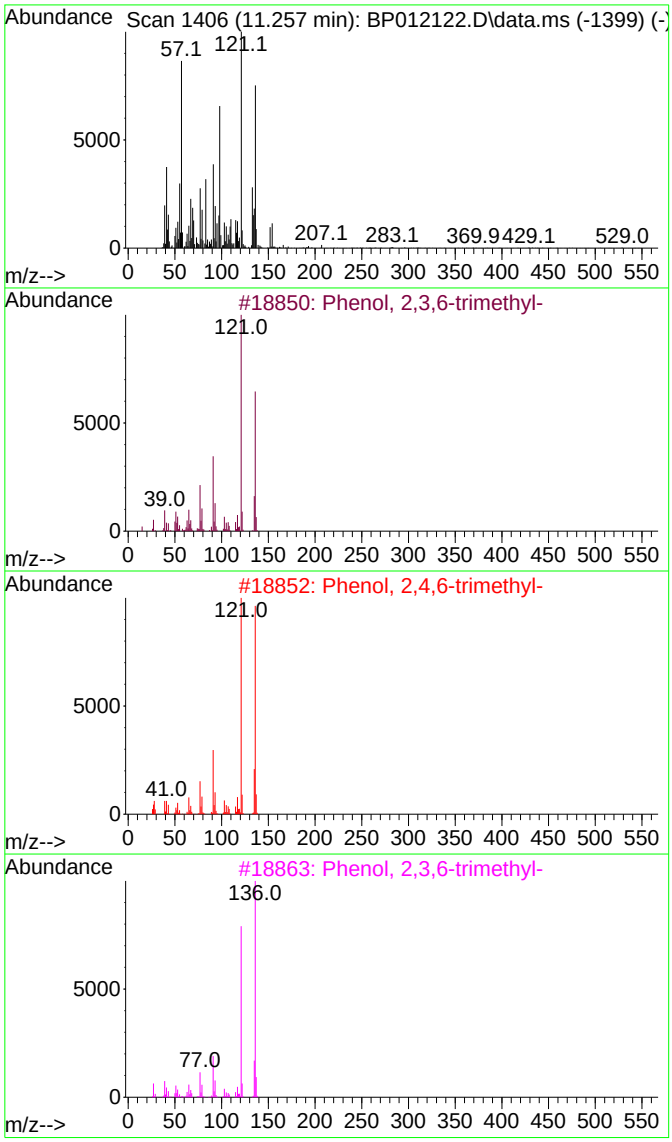
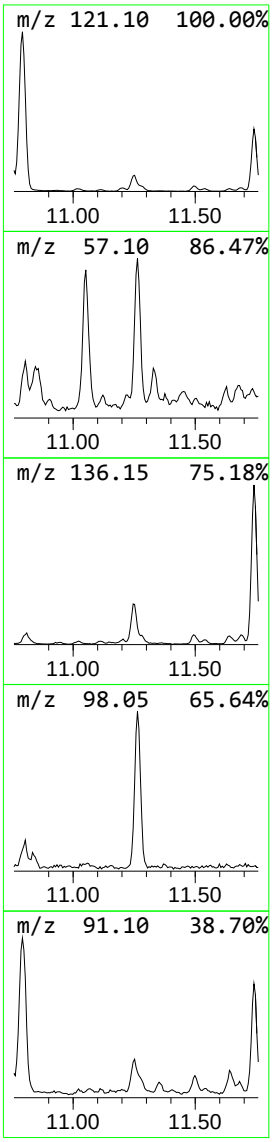
Peak Number 22 Phenol, 2,3,6-trimethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.258	3.27 ng/ul	304438	Naphthalene-d8	10.658

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	91
2	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	83
3	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	83
4	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	64
5	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	60





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ClientSampleId :
BHA38

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

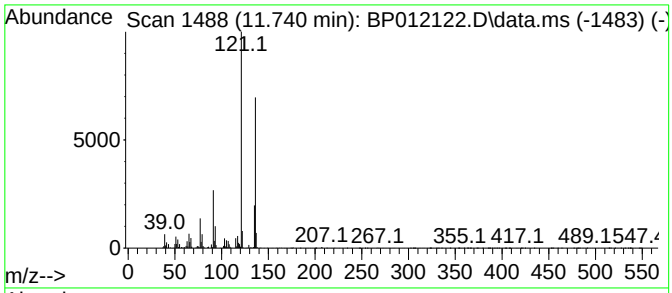
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

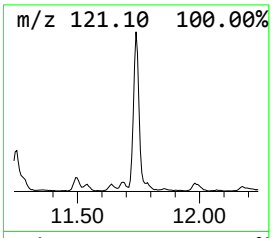
****

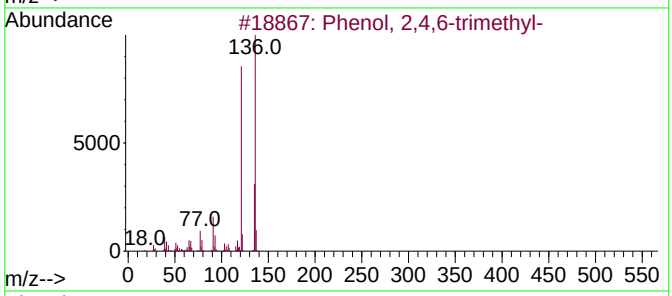
Peak Number 23 Phenol, 2,4,6-trimethyl- Concentration Rank 15

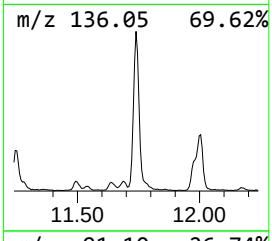
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.740	4.44 ng/ul	412859	Naphthalene-d8	10.658

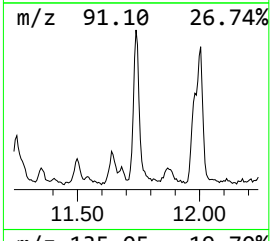
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	97
2	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	96
3	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	96
4	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	95
5	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	95

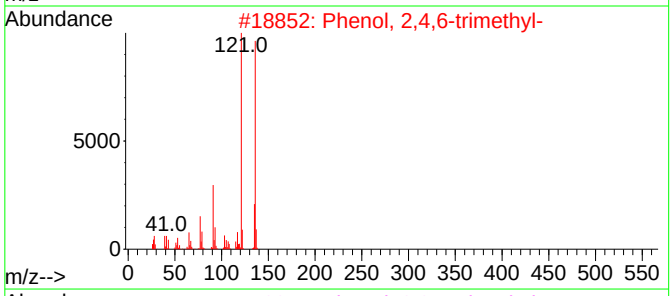


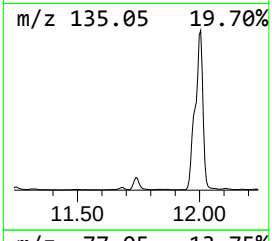


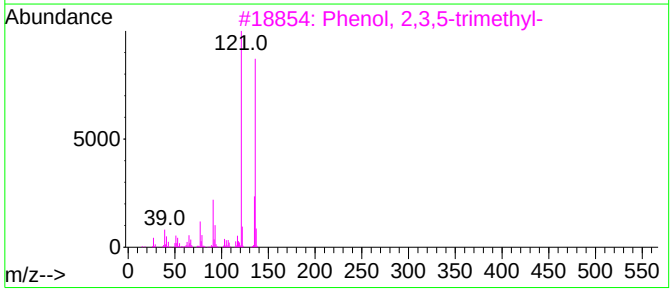


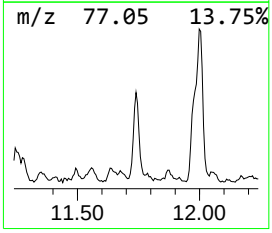












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Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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

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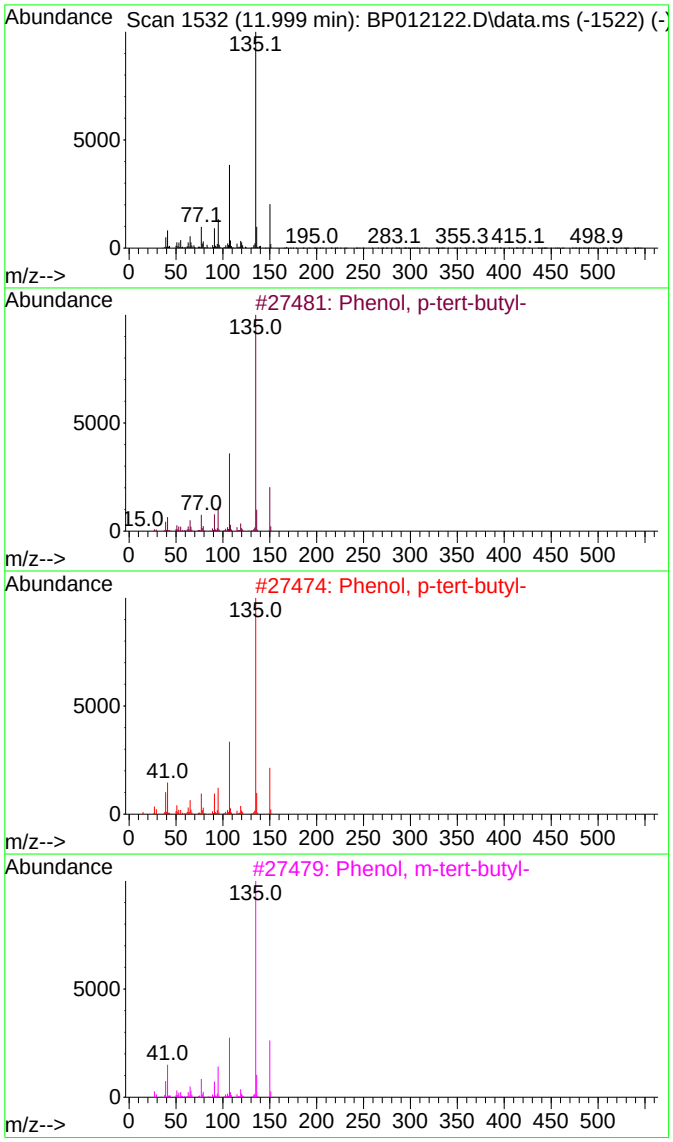
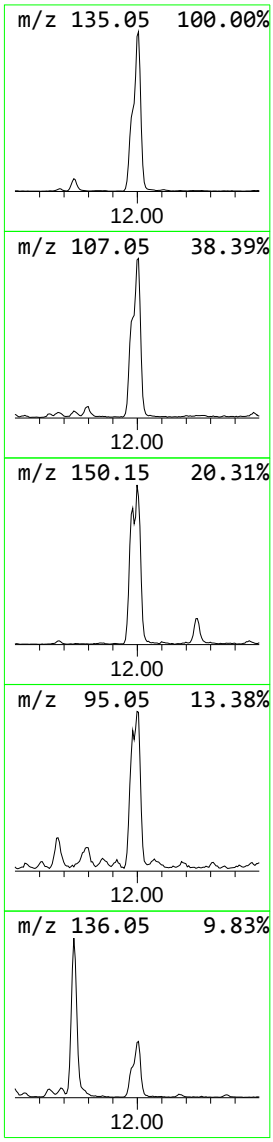
Peak Number 24 Phenol, p-tert-butyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.999	14.76 ng/ul	1372370	Naphthalene-d8	10.658

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	94
5	4-Hydroxy-2-methylacetophenone	150	C9H10O2	000875-59-2	87





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Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

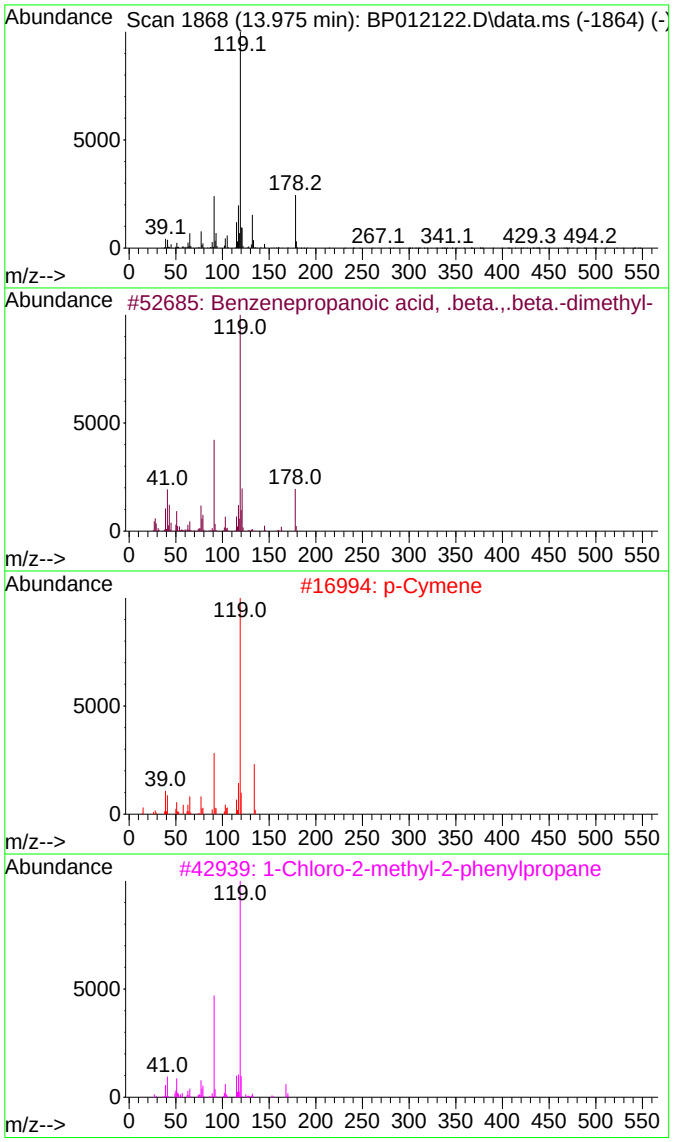
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

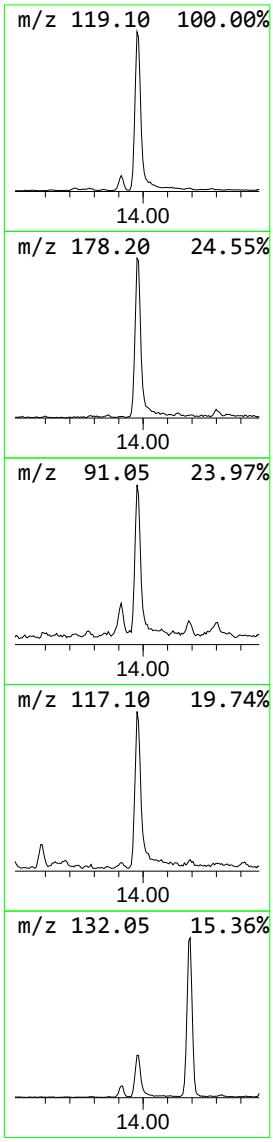
****

Peak Number 25 Benzenepropanoic acid, .bet... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.975	3.00 ng/ul	420933	Acenaphthene-d10	14.499

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenepropanoic acid, .beta.,.b...	178	C11H14O2	001010-48-6	64
2	p-Cymene	134	C10H14	000099-87-6	50
3	1-Chloro-2-methyl-2-phenylpropane	168	C10H13Cl	000515-40-2	50
4	o-Cymene	134	C10H14	000527-84-4	49
5	1-Piperidinecarbothioic acid, S-...	263	C15H21NOS	061432-55-1	47





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Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

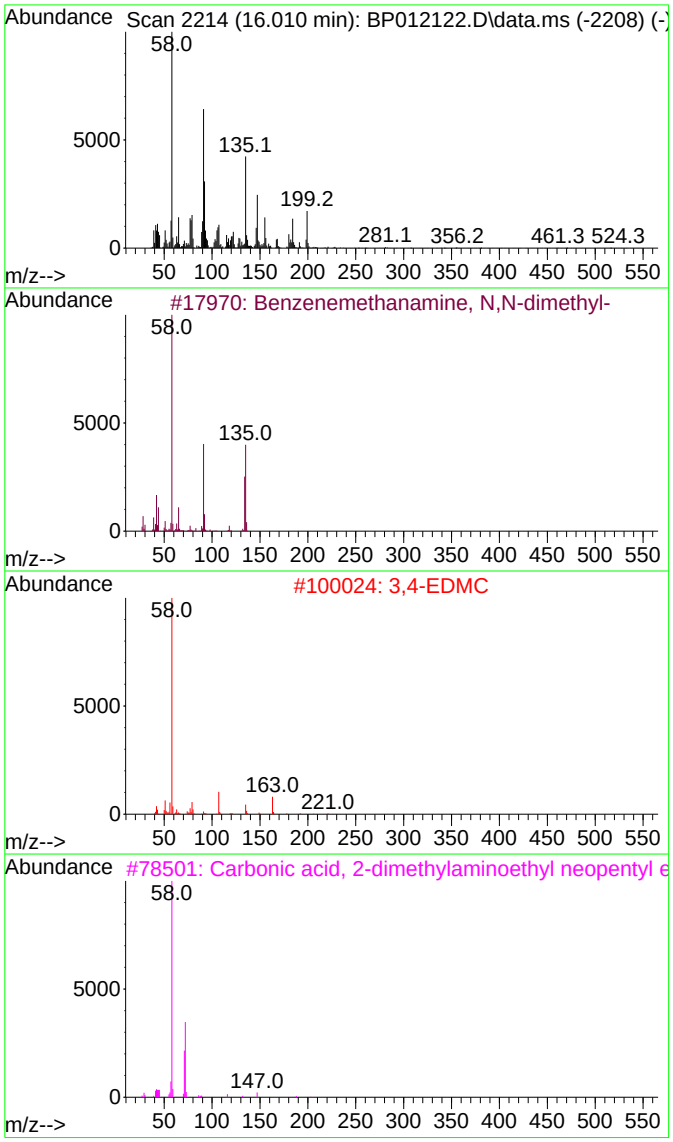
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

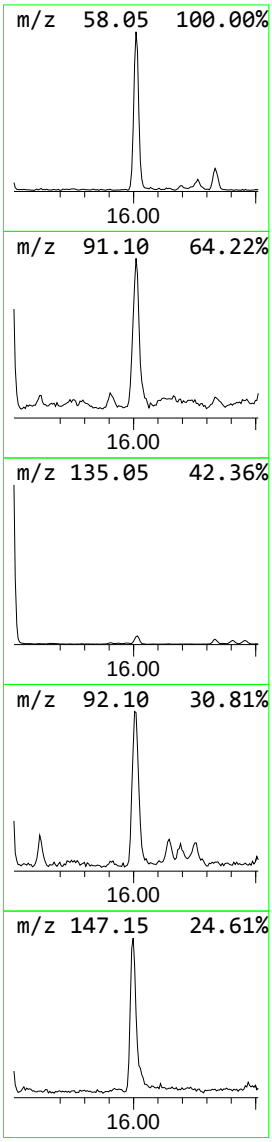
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 Benzenemethanamine, N,N-dim... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.010	3.45 ng/ul	451990	Phenanthrene-d10	17.257

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenemethanamine, N,N-dimethyl-	135	C9H13N	000103-83-3	53
2	3,4-EDMC	221	C12H15NO3	030253-44-2	30
3	Carbonic acid, 2-dimethylaminoet...	203	C10H21NO3	1000331-43-1	30
4	1-(4-Amino-furazan-3-yl)-5-dimet...	253	C8H11N7O3	1000276-04-0	27
5	Amine, dimethyl-2-phosphinoethyl-	105	C4H12NP	161944-90-7	25





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Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

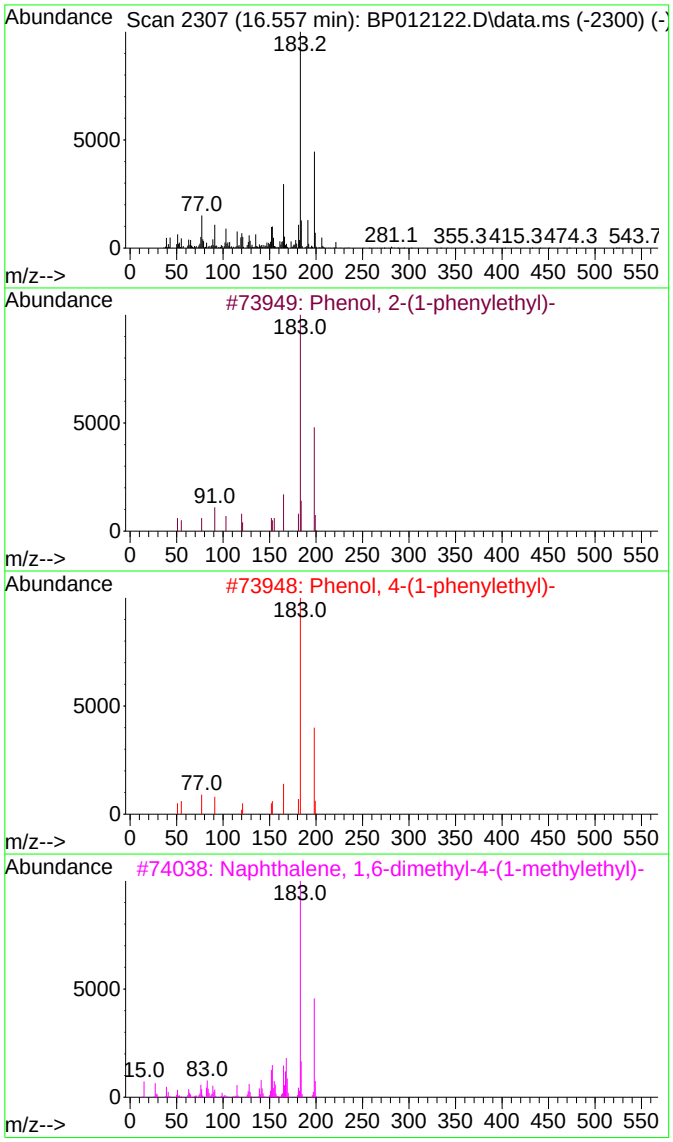
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

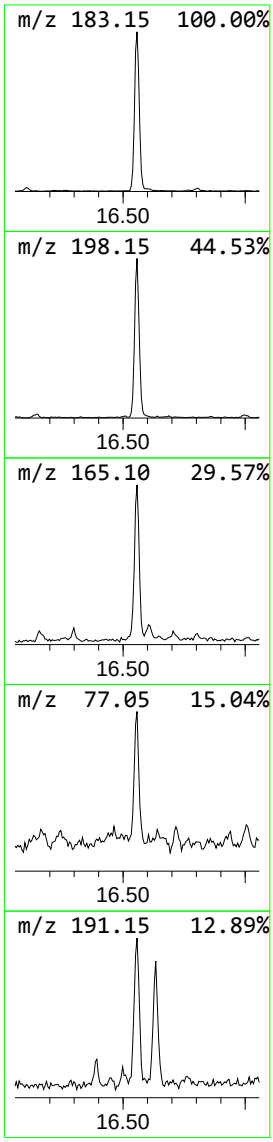
****

Peak Number 27 Phenol, 2-(1-phenylethyl)- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.557	2.77 ng/ul	363147	Phenanthrene-d10	17.257

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1-phenylethyl)-	198	C14H14O	004237-44-9	91
2	Phenol, 4-(1-phenylethyl)-	198	C14H14O	001988-89-2	81
3	Naphthalene, 1,6-dimethyl-4-(1-m...	198	C15H18	000483-78-3	74
4	Azulene, 1,4-dimethyl-7-(1-methy...	198	C15H18	000489-84-9	74
5	Naphthalene, 1,6-dimethyl-4-(1-m...	198	C15H18	000483-78-3	72





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Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

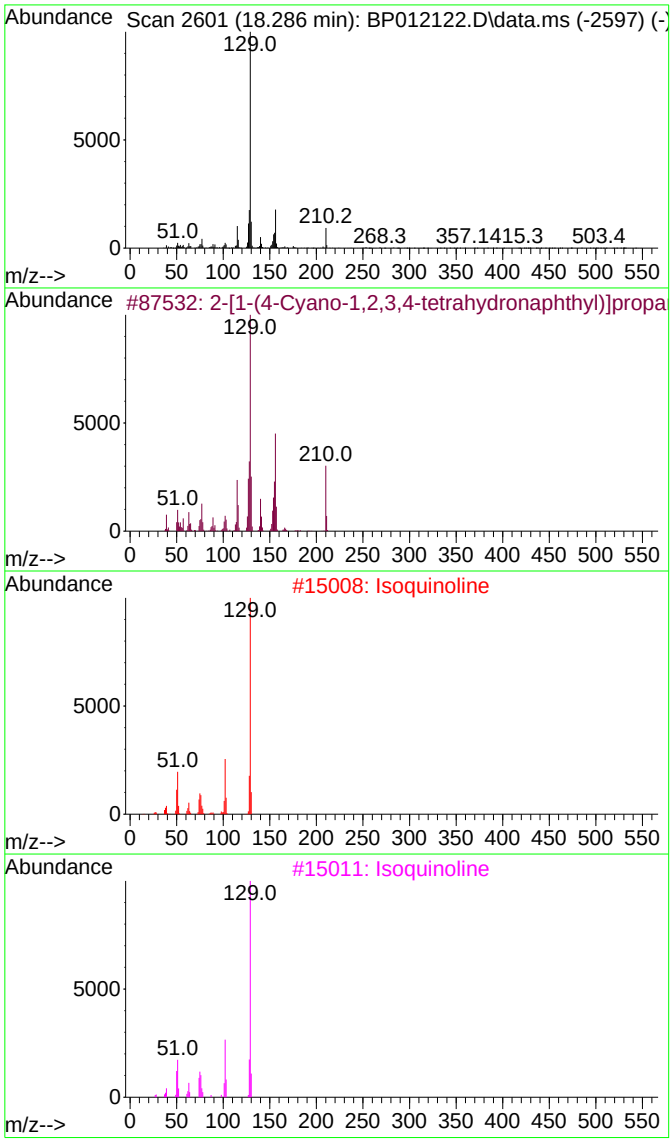
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

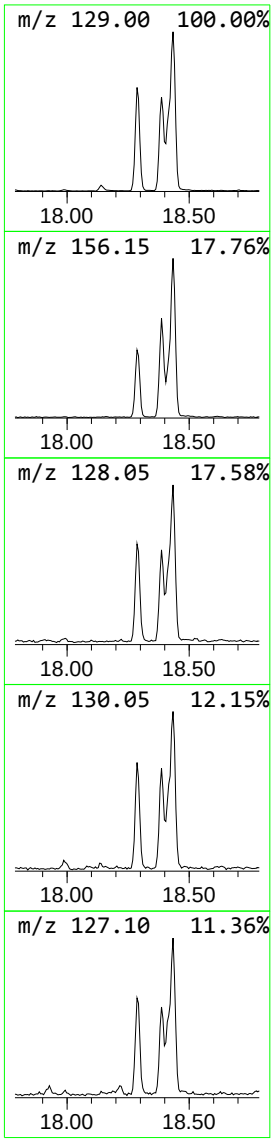
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 2-[1-(4-Cyano-1,2,3,4-tetra... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.287	2.70 ng/ul	353776	Phenanthrene-d10	17.257

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-[1-(4-Cyano-1,2,3,4-tetrahydro...	210	C14H14N2	057964-39-3	78
2	Isoquinoline	129	C9H7N	000119-65-3	52
3	Isoquinoline	129	C9H7N	000119-65-3	52
4	1-Deuteronaphthalene	129	C10H7D	000875-62-7	50
5	Quinoline-8-carboxylic acid	173	C10H7NO2	000086-59-9	42





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Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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

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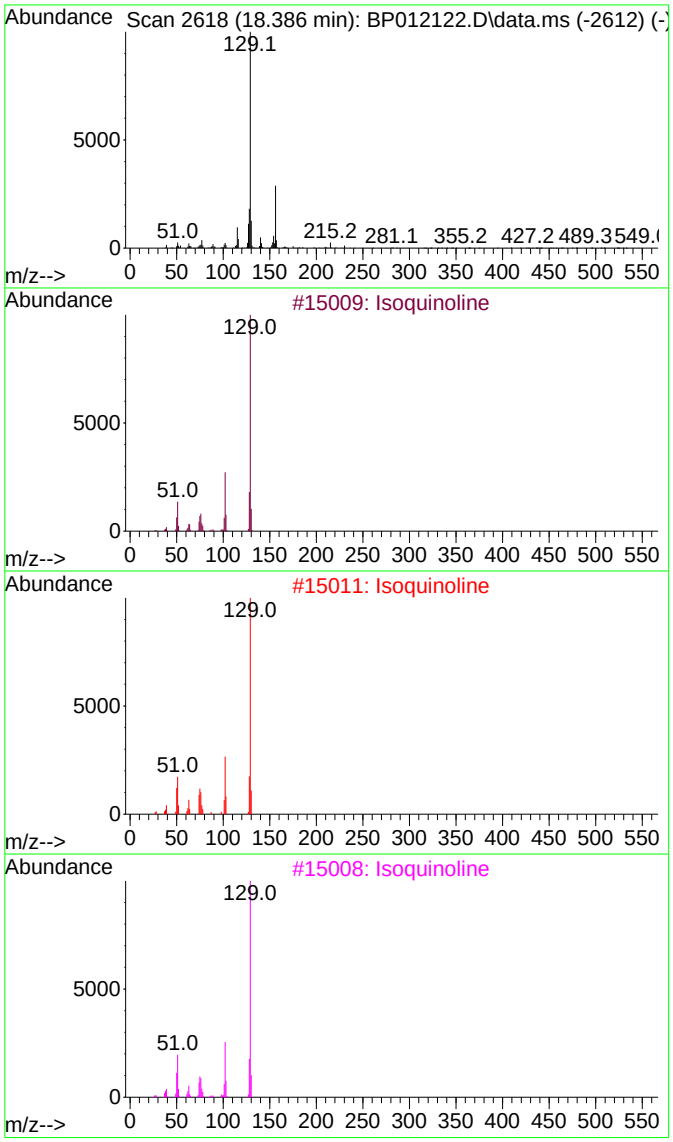
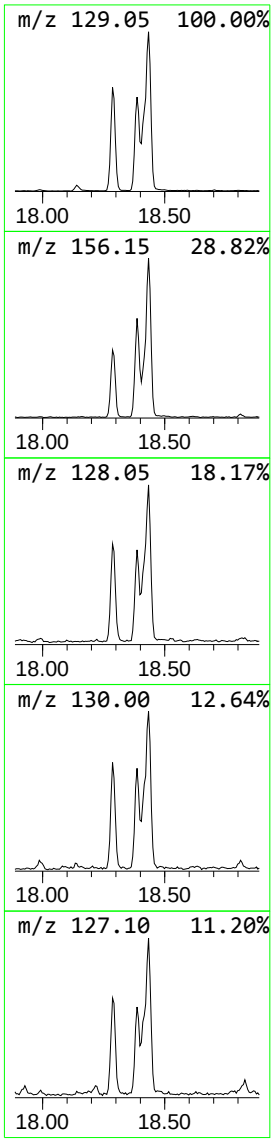
Peak Number 29 Isoquinoline Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.387	2.48 ng/ul	324711	Phenanthrene-d10	17.257

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isoquinoline	129	C9H7N	000119-65-3	52
2	Isoquinoline	129	C9H7N	000119-65-3	50
3	Isoquinoline	129	C9H7N	000119-65-3	50
4	1,2-Benzenediacetonitrile	156	C10H8N2	000613-73-0	50
5	3-[1-(4-Cyano-1,2,3,4-tetrahydro...	210	C14H14N2	057964-40-6	42





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

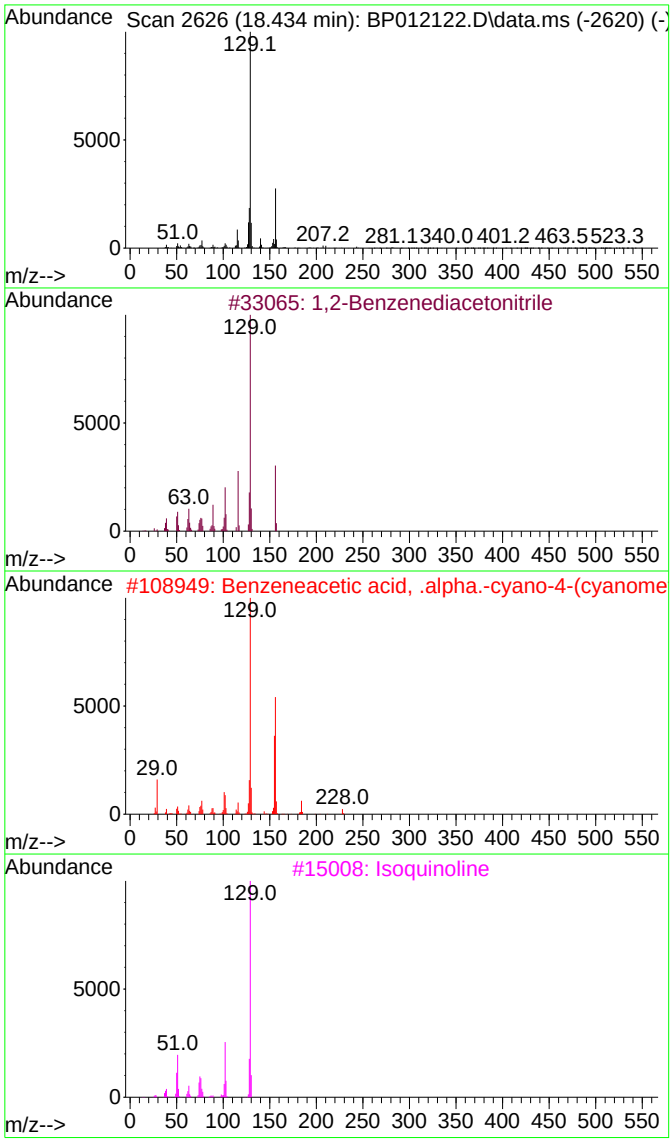
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

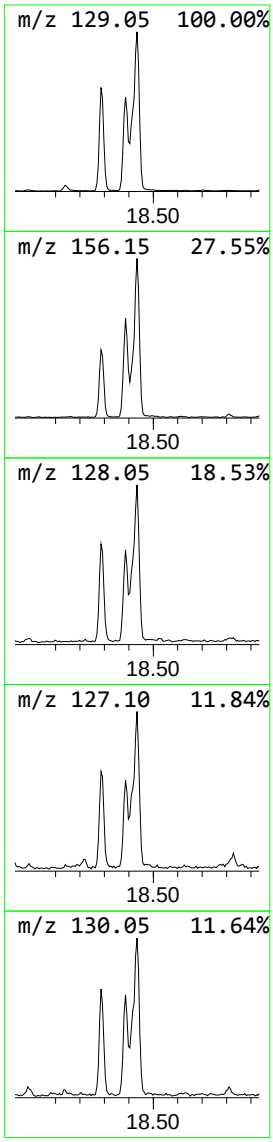
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 30 1,2-Benzenediacetonitrile Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.434	5.35 ng/ul	701960	Phenanthrene-d10	17.257

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenediacetonitrile	156	C10H8N2	000613-73-0	64
2	Benzeneacetic acid, .alpha.-cyan...	228	C13H12N2O2	134882-04-5	56
3	Isoquinoline	129	C9H7N	000119-65-3	52
4	Quinoline	129	C9H7N	000091-22-5	50
5	Isoquinoline	129	C9H7N	000119-65-3	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
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Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

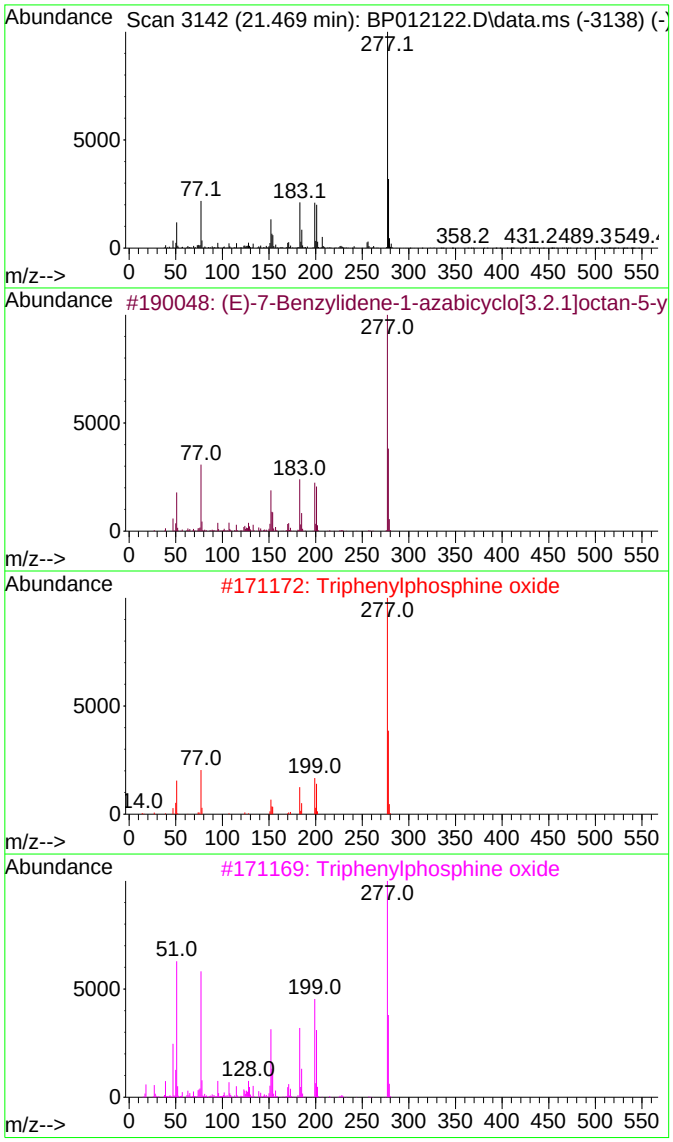
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

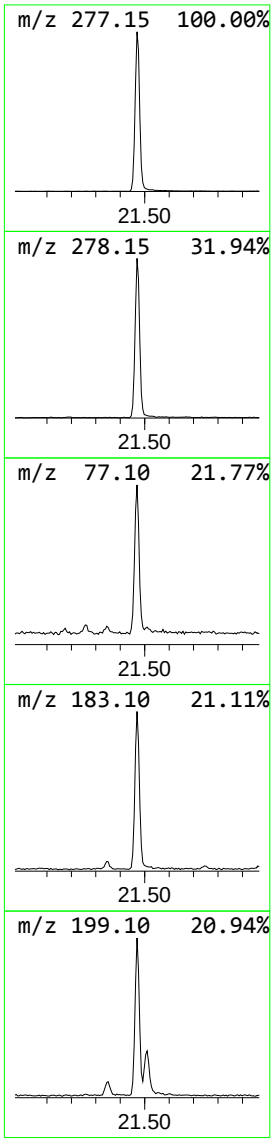
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 31 (E)-7-Benzylidene-1-azabicy... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.469	4.14 ng/ul	689431	Chrysene-d12	21.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-7-Benzylidene-1-azabicyclo[3...	293	C15H19NO3S	1000483-83-4	91
2	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	81
3	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	78
4	Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	64
5	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012122.D
Acq On : 13 Oct 2022 15:54
Operator : CG/JU
Sample : N5013-01
Misc :
ALS Vial : 93 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA38

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.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
Triethylamine	3.070	12.1	ng/ul	853287	1	7.846	1412710	20.0
2-Pentanol, 4-m...	3.840	2.2	ng/ul	154976	1	7.846	1412710	20.0
Benzene, 1,3-di...	5.475	75.2	ng/ul	5310470	1	7.846	1412710	20.0
3-Heptanone, 6-...	6.511	3.1	ng/ul	216539	1	7.846	1412710	20.0
Benzene, 1-ethy...	6.928	4.9	ng/ul	346367	1	7.846	1412710	20.0
Dill ether	7.552	2.3	ng/ul	165366	1	7.846	1412710	20.0
Furan, tetrahyd...	8.046	3.3	ng/ul	231638	1	7.846	1412710	20.0
Indane	8.216	14.7	ng/ul	1037410	1	7.846	1412710	20.0
Cyclohexanol, 3...	8.475	45.0	ng/ul	3174900	1	7.846	1412710	20.0
Aniline, N-methyl-	8.852	2.1	ng/ul	147883	1	7.846	1412710	20.0
Benzene, 1-ethy...	8.952	3.5	ng/ul	249803	1	7.846	1412710	20.0
Benzene, 1-(chl...	9.081	3.2	ng/ul	226115	1	7.846	1412710	20.0
3-Octanol, 3,7-...	9.169	7.5	ng/ul	530738	1	7.846	1412710	20.0
(+)-2-Bornanone	10.069	5.5	ng/ul	508052	2	10.658	1860070	20.0
Bicyclo[4.1.0]h...	11.069	4.1	ng/ul	382578	2	10.658	1860070	20.0
Phenol, 2,3,6-t...	11.258	3.3	ng/ul	304438	2	10.658	1860070	20.0
Phenol, 2,4,6-t...	11.740	4.4	ng/ul	412859	2	10.658	1860070	20.0
Phenol, p-tert-...	11.999	14.8	ng/ul	1372370	2	10.658	1860070	20.0
Benzenepropanoi...	13.975	3.0	ng/ul	420933	3	14.499	2805870	20.0
Benzenemethanam...	16.010	3.5	ng/ul	451990	4	17.257	2622050	20.0
Phenol, 2-(1-ph...	16.557	2.8	ng/ul	363147	4	17.257	2622050	20.0
2-[1-(4-Cyano-1...	18.287	2.7	ng/ul	353776	4	17.257	2622050	20.0
Isoquinoline	18.387	2.5	ng/ul	324711	4	17.257	2622050	20.0
1,2-Benzenediac...	18.434	5.3	ng/ul	701960	4	17.257	2622050	20.0
(E)-7-Benzylide...	21.469	4.1	ng/ul	689431	5	21.345	3331650	20.0