LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012812.D
Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0L03

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-BP112522.M
Title : SVOA CALIBRATION

Signal : TIC: BP012812.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.393	32	35	40	rBV	84065	116744	0.69%	0.055%
2	3.822	106	108	122	rBV	446822	738833	4.37%	0.348%
3	4.058	144	148	158	rVB	52279	89608	0.53%	0.042%
4	4.269	178	184	196	rVV2	75556	158492	0.94%	0.075%
5	6.075	482	491	501	rBV	192296	338063	2.00%	0.159%
6	6.652	582	589	597	rBV	242935	369364	2.18%	0.174%
7	7.163	669	676	686	rVB2	475064	990824	5.86%	0.467%
8	7.340	699	706	719	rVB	1895687	2961169	17.51%	1.395%
9	7.475	722	729	734	rBV	91366	221850	1.31%	0.105%
10	7.546	734	741	748	rVV	1644182	2707968	16.01%	1.276%
11	7.610	748	752	761	rVB2	59856	112167	0.66%	0.053%
12	7.793	777	783	793	rVB2	127667	242219	1.43%	0.114%
13	8.016	812	821	827	rBV	1929520	3063392	18.11%	1.443%
14	8.075	827	831	837	rVB	95479	149328	0.88%	0.070%
15	8.716	932	940	950	rVV	1452059	2325760	13.75%	1.096%
16	8.810	950	956	957	rVV	67421	104516	0.62%	0.049%
17	8.840	957	961	967	rVB	152630	256733	1.52%	0.121%
18	9.122	1002	1009	1013	rBV3	61972	113269	0.67%	0.053%
19	9.181	1013	1019	1028	rVV	2027416	3282019	19.40%	1.546%
20	9.263	1028	1033	1040	rVB3	99801	192309	1.14%	0.091%
21	9.346	1040	1047	1056	rBV	307130	497844	2.94%	0.235%
22	9.904	1134	1142	1157	rBV	1391945	2395667	14.16%	1.129%
23	10.075	1157	1171	1175	rVV3	118081	415221	2.45%	0.196%
24	10.122	1175	1179	1189	rVB2	51554	111182	0.66%	0.052%
25	10.310	1204	1211	1215	rBV	179789	306629	1.81%	0.144%
26	10.446	1227	1234	1243	rBV	2596935	4223600	24.97%	1.990%
27	10.604	1254	1261	1269	rBV2	534530	937623	5.54%	0.442%
28	10.693	1270	1276	1281	rVV4	144259	270055	1.60%	0.127%
29	10.834	1291	1300	1307	rBV	2927732	4859896	28.73%	2.290%
30	10.963	1313	1322	1334	rBV	2279348	3904624	23.08%	1.840%
31	11.187	1354	1360	1373	rVB	330025	664286	3.93%	0.313%
32	11.345	1381	1387	1395	rBV2	212187	465947	2.75%	0.220%
33	11.422	1395	1400	1405	rBV	111257	174956	1.03%	0.082%
34	11.481	1405	1410	1420	rVB	337545	595833	3.52%	0.281%
35	11.569	1420	1425	1430	rBV2	103006	187116	1.11%	0.088%
36	11.622	1430	1434	1444	rVB6	45633	109434	0.65%	0.052%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012812.D
Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

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-BP112522.M
Title : SVOA CALIBRATION

37	11.722	1444	1451	1455	rBV2	64993	119803	0.71%	0.056%
38	11.851	1467	1473	1483	rVB10	31213	90726	0.54%	0.043%
39	12.081	1504	1512	1516	rBV2	176728	328351	1.94%	0.155%
40	12.287	1542	1547	1549	rBV	61725	98012	0.58%	0.046%
41	12.334	1550	1555	1560	rVB3	110750	221021	1.31%	0.104%
42	12.410	1565	1568	1576	rVB3	57302	92349	0.55%	0.044%
43	12.563	1587	1594	1599	rVB3	106931	209186	1.24%	0.099%
44	12.845	1634	1642	1646	rBV2	162419	301991	1.79%	0.142%
45	12.887	1646	1649	1654	rVV	156287	232166	1.37%	0.109%
46	12.951	1654	1660	1664	rVB5	52652	98506	0.58%	0.046%
47	13.016	1664	1671	1683	rBV	6630045	10347966	61.17%	4.876%
48	13.104	1683	1686	1692	rVB	236527	338168	2.00%	0.159%
49	13.263	1705	1713	1727	rBV	9962976	14551748	86.02%	6.856%
50	13.398	1732	1736	1740	rVB	86005	110674	0.65%	0.052%
51	13.492	1746	1752	1757	rBV	2797594	3930830	23.24%	1.852%
52	13.545	1757	1761	1769	rVB	638676	980397	5.80%	0.462%
53	13.957	1827	1831	1836	rVB	337989	464874	2.75%	0.219%
54	14.057	1842	1848	1855	rBV	5573751	7754179	45.84%	3.653%
55	14.181	1862	1869	1873	rVB4	91437	164826	0.97%	0.078%
56	14.298	1883	1889	1891	rBV3	145708	235359	1.39%	0.111%
57	14.345	1891	1897	1903	rVV	6186086	9070080	53.62%	4.273%
58	14.398	1903	1906	1917	rVB2	167180	310472	1.84%	0.146%
59	14.651	1936	1949	1956	rBV	4917398	7266646	42.96%	3.424%
60	14.834	1971	1980	1989	rBV	356938	611633	3.62%	0.288%
61	14.922	1990	1995	1999	rVB2	59269	91356	0.54%	0.043%
62	15.057	2014	2018	2027	rVB7	43407	93271	0.55%	0.044%
63	15.239	2045	2049	2055	rVB2	142337	208357	1.23%	0.098%
64	15.345	2063	2067	2069	rBV	61816	90717	0.54%	0.043%
65	15.386	2069	2074	2081	rVV	1672934	2216980	13.11%	1.045%
66	15.463	2081	2087	2095	rVB	2831824	3839305	22.70%	1.809%
67	15.645	2111	2118	2130	rVB	8777690	11859341	70.11%	5.588%
68	15.751	2130	2136	2145	rBV	1675421	2182073	12.90%	1.028%
69	16.004	2172	2179	2187	rBV2	1275967	1815646	10.73%	0.855%
70	16.098	2192	2195	2200	rVB4	62722	88998	0.53%	0.042%
71	16.257	2218	2222	2225	rBV	102067	140788	0.83%	0.066%
72	16.292	2225	2228	2233	rVV	98859	125326	0.74%	0.059%
73	16.345	2233	2237	2242	rVB	159716	216939	1.28%	0.102%
74	16.439	2250	2253	2257	rVB	83513	102616	0.61%	0.048%
75	16.651	2284	2289	2295	rBV	403406	560315	3.31%	0.264%
76	16.763	2301	2308	2315	rBV3	204247	452944	2.68%	0.213%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012812.D
Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

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-BP112522.M
Title : SVOA CALIBRATION

77	16.857	2320	2324	2328	rBV2	59809	98281	0.58%	0.046%
78	17.157	2372	2375	2379	rVV	243060	279582	1.65%	0.132%
79	17.198	2379	2382	2388	rVB2	120685	167903	0.99%	0.079%
80	17.263	2389	2393	2403	rVB2	93116	144954	0.86%	0.068%
81	17.357	2404	2409	2412	rBV5	66646	108983	0.64%	0.051%
82	17.404	2412	2417	2427	rVB	5955753	8642316	51.09%	4.072%
83	17.504	2428	2434	2444	rBV	9354717	13633513	80.60%	6.424%
84	17.686	2462	2465	2474	rVV	195726	285096	1.69%	0.134%
85	17.763	2475	2478	2485	rVB2	54786	88598	0.52%	0.042%
86	17.969	2509	2513	2518	rBV	60583	83633	0.49%	0.039%
87	18.069	2525	2530	2534	rBV2	213277	261851	1.55%	0.123%
88	18.269	2558	2564	2568	rVB5	155700	294416	1.74%	0.139%
89	18.363	2573	2580	2596	rVB	9373051	13139101	77.67%	6.191%
90	18.645	2624	2628	2632	rBV3	100853	128203	0.76%	0.060%
91	18.775	2648	2650	2659	rVV5	66762	110019	0.65%	0.052%
92	18.845	2659	2662	2669	rVB2	91415	126964	0.75%	0.060%
93	19.310	2737	2741	2748	rVB	124960	167972	0.99%	0.079%
94	19.374	2748	2752	2757	rBV	130917	193711	1.15%	0.091%
95	19.439	2758	2763	2769	rVV5	124837	184324	1.09%	0.087%
96	19.504	2771	2774	2778	rVV2	81997	92585	0.55%	0.044%
97	19.733	2807	2813	2822	rBV	12980120	16915857	100.00%	7.970%
98	21.492	3106	3112	3124	rBV	7420802	9985975	59.03%	4.705%
99	23.845	3504	3512	3520	rBV2	8053516	16555060	97.87%	7.800%
100	24.010	3532	3540	3557	rVB2	4826897	9882321	58.42%	4.656%

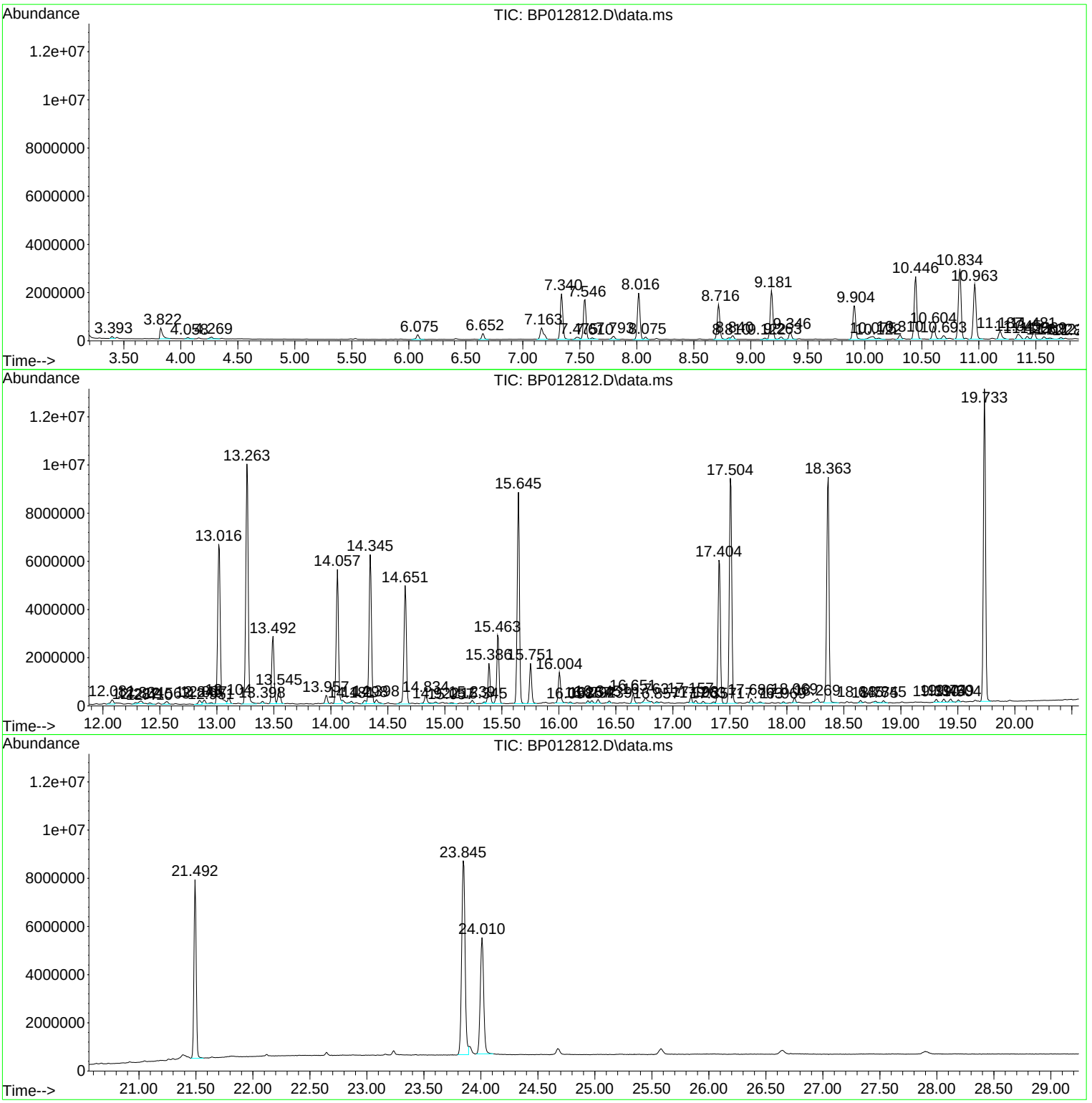
Sum of corrected areas: 212240693

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012812.D
Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012812.D
Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

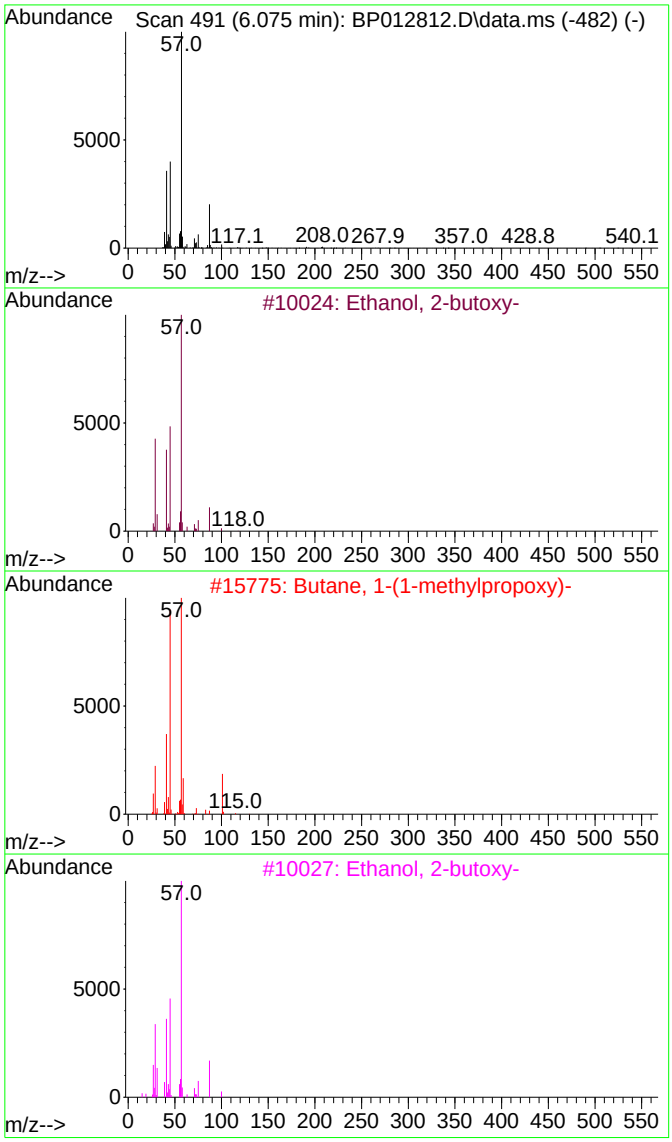
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

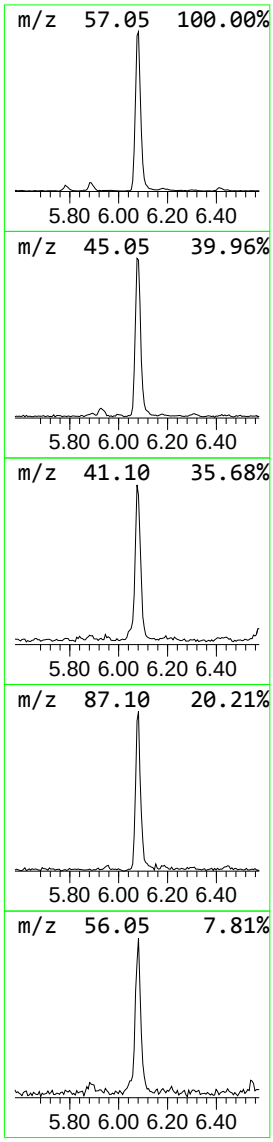
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Ethanol, 2-butoxy- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.075	2.21 ng/ul	338063	1,4-Dichlorobenzene-d4	8.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	50
2	Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	50
3	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	47
4	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	47
5	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012812.D
Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

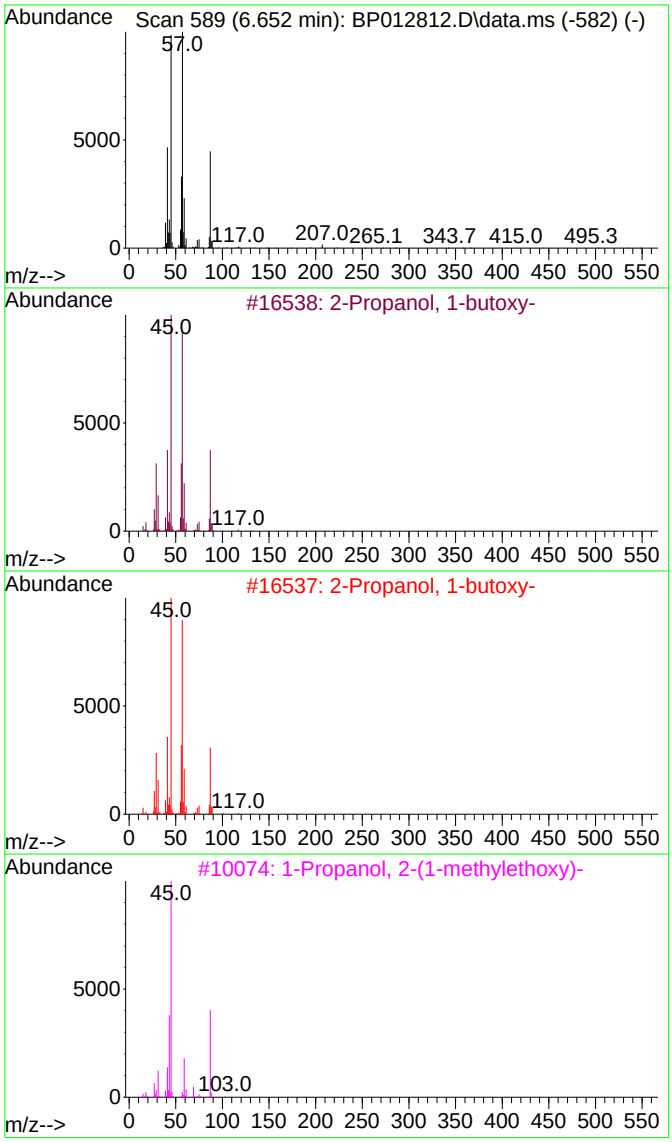
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

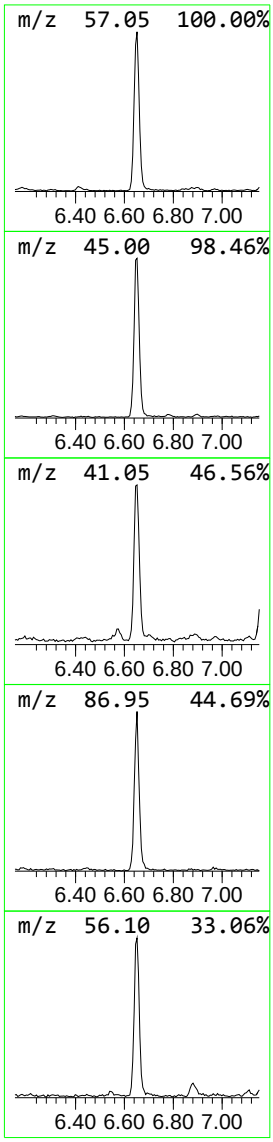
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Propanol, 1-butoxy- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.652	2.41 ng/ul	369364	1,4-Dichlorobenzene-d4	8.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-butoxy-	132	C7H16O2	005131-66-8	90
2	2-Propanol, 1-butoxy-	132	C7H16O2	005131-66-8	72
3	1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	59
4	1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	59
5	Carbonic acid, bis(1-methylpropy...	174	C9H18O3	000623-63-2	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012812.D
Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

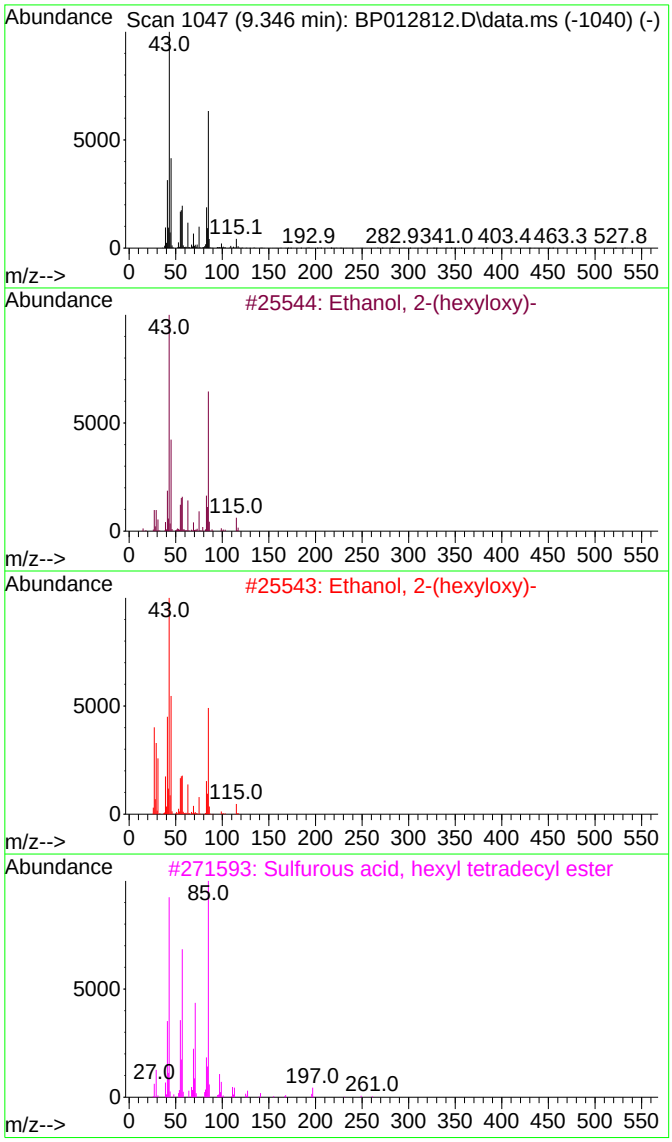
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

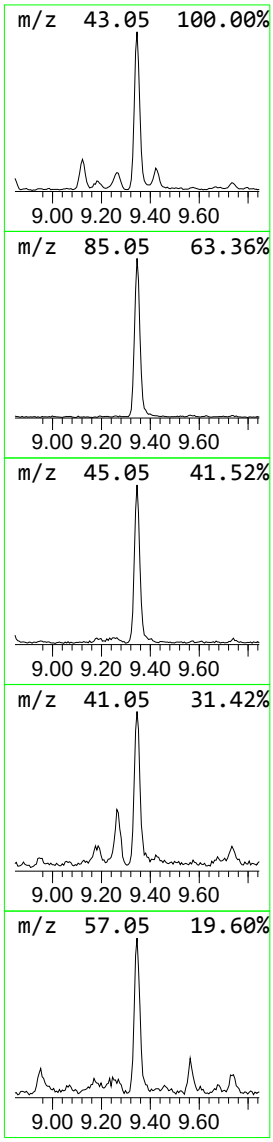
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Ethanol, 2-(hexyloxy)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.346	3.25 ng/ul	497844	1,4-Dichlorobenzene-d4	8.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	83
2	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	64
3	Sulfurous acid, hexyl tetradecyl...	362	C20H42O3S	1010309-13-6	47
4	Hexanal dihexyl acetal	286	C18H38O2	1000430-99-9	43
5	2-Butanol, 1,4-dimethoxy-3-(1,2,...	234	C12H26O4	079449-92-6	40





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012812.D
Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

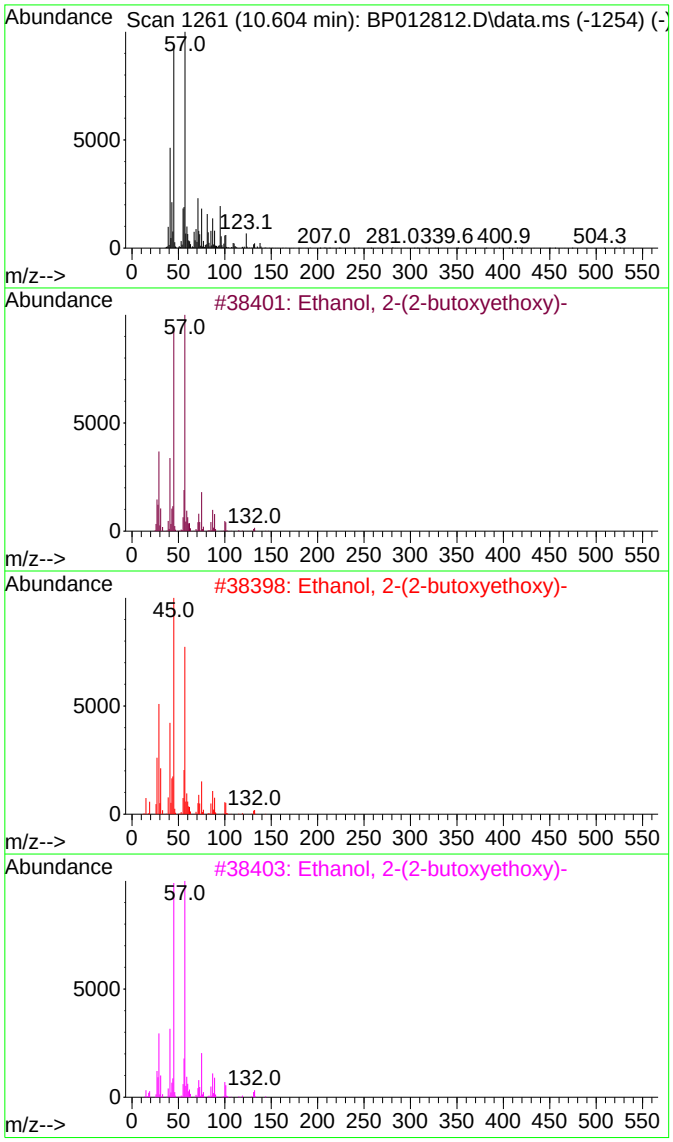
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

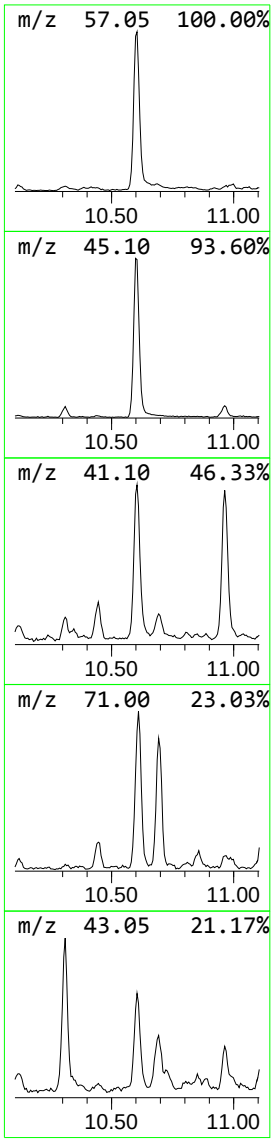
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.604	3.86 ng/ul	937623	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	58
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	58
3	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	58
4	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	53
5	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012812.D
Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

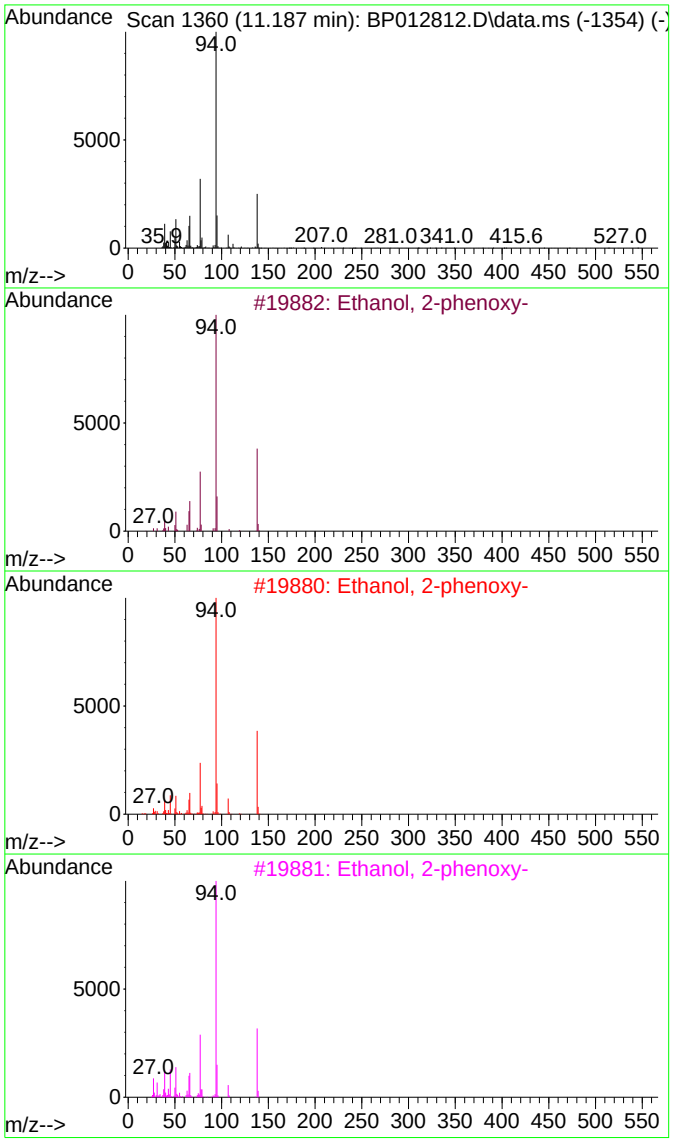
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

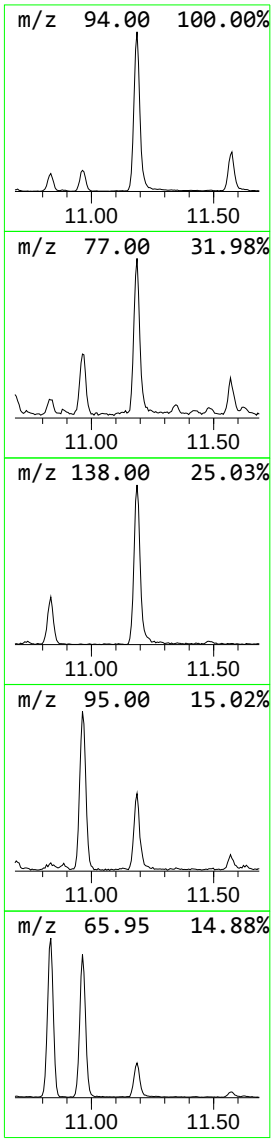
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Ethanol, 2-phenoxy- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.187	2.73 ng/ul	664286	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	95
2	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	94
3	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	91
4	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	64
5	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	64





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Data File : BP012812.D
Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

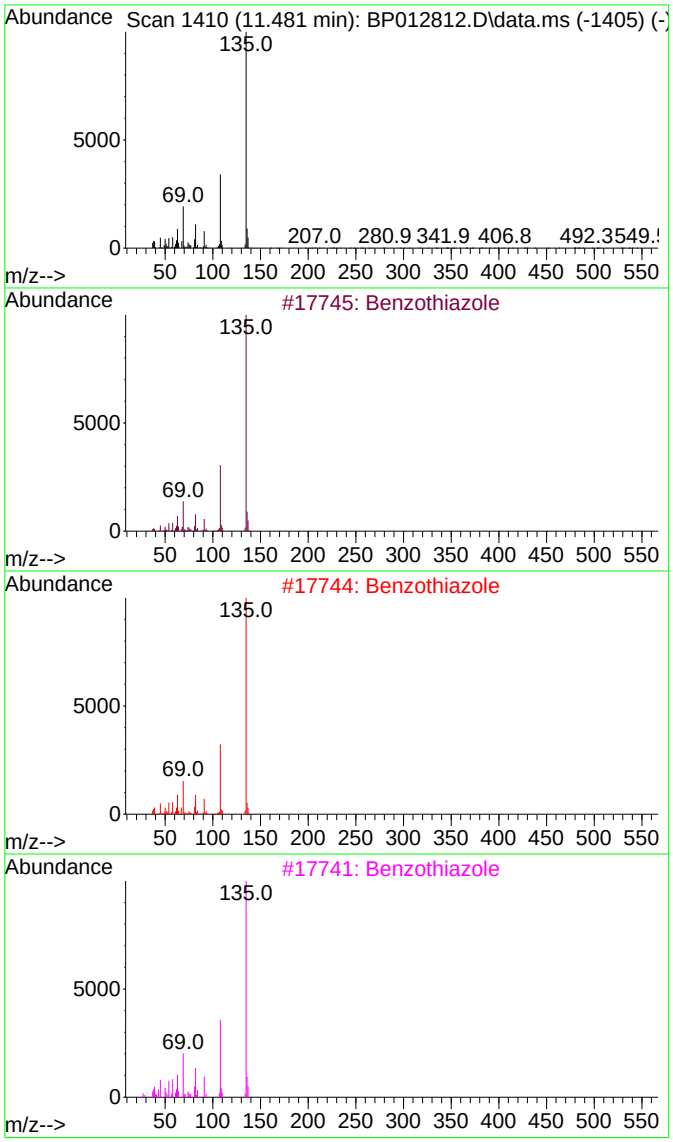
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

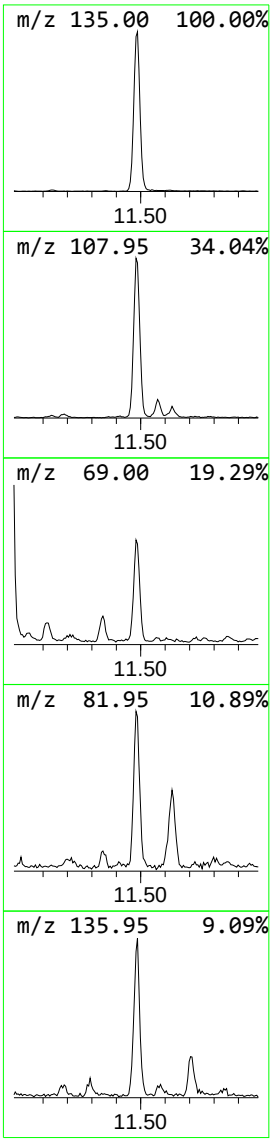
****

Peak Number 6 Benzothiazole Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.481	2.45 ng/ul	595833	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzothiazole	135	C7H5NS	000095-16-9	95
2	Benzothiazole	135	C7H5NS	000095-16-9	95
3	Benzothiazole	135	C7H5NS	000095-16-9	95
4	Benzothiazole	135	C7H5NS	000095-16-9	91
5	Benzothiazole	135	C7H5NS	000095-16-9	91





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Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

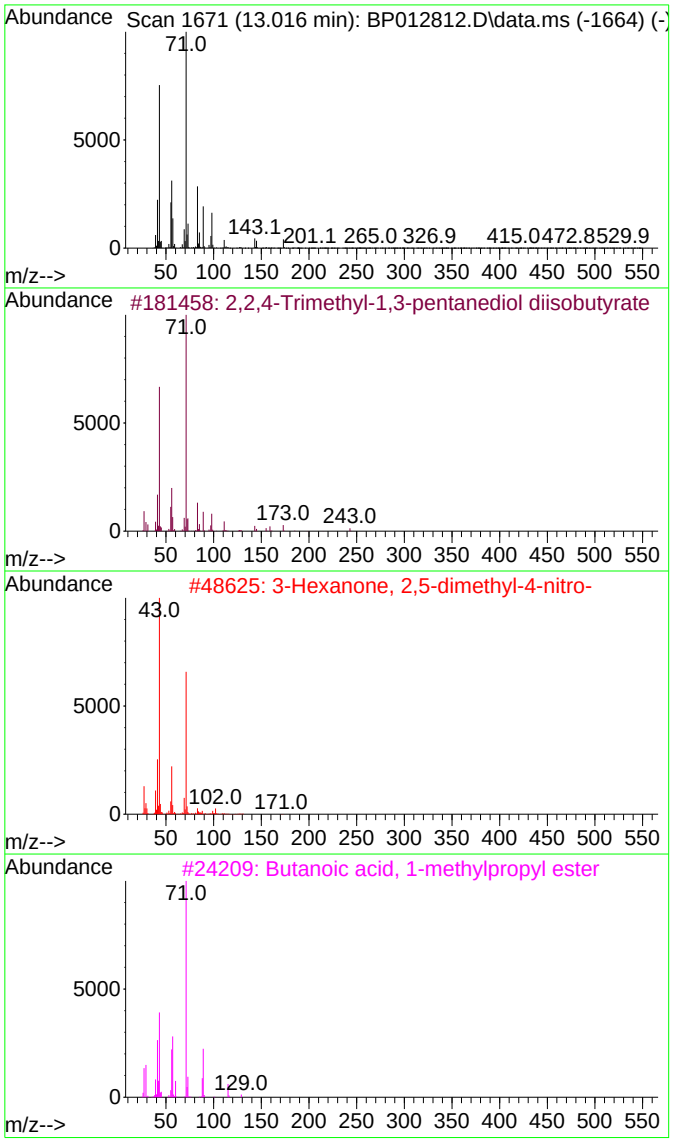
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

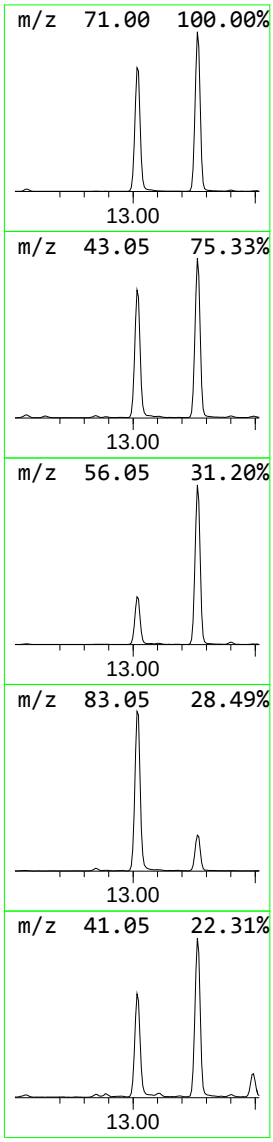
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 2,2,4-Trimethyl-1,3-pentane... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.016	28.48 ng/ul	10348000	Acenaphthene-d10	14.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	78
2	3-Hexanone, 2,5-dimethyl-4-nitro-	173	C8H15NO3	059906-54-6	43
3	Butanoic acid, 1-methylpropyl ester	144	C8H16O2	000819-97-6	42
4	Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	40
5	Isophytol	296	C20H40O	000505-32-8	38





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Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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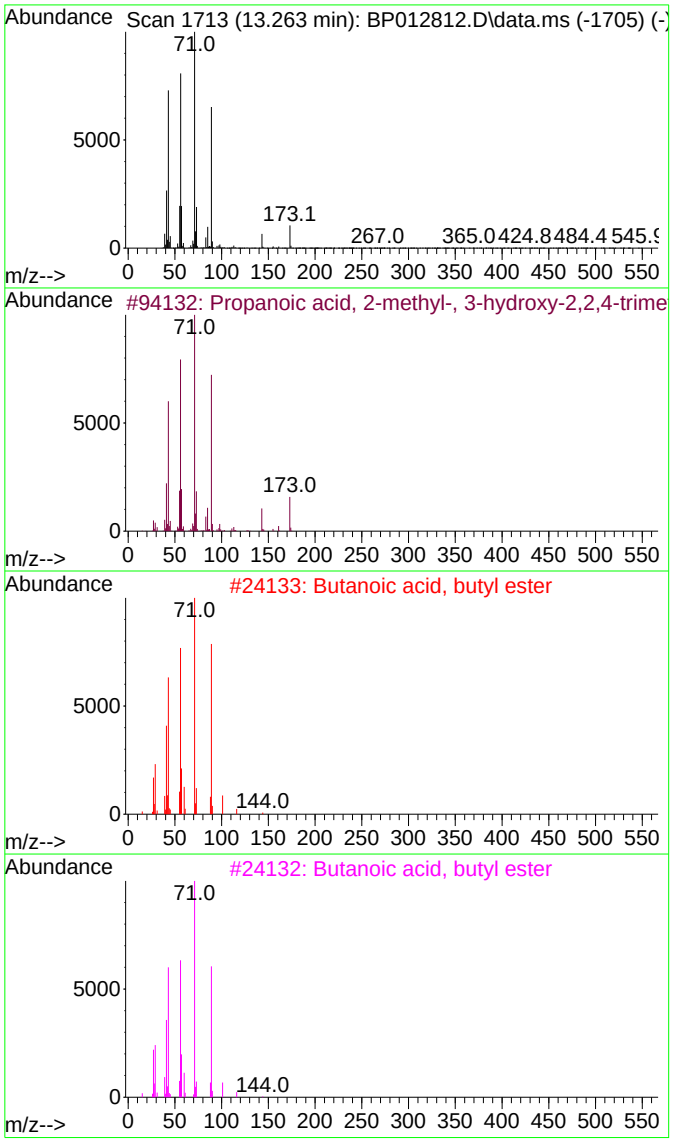
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

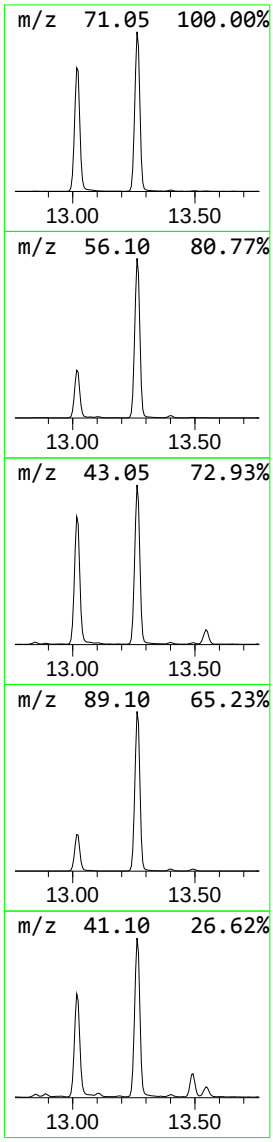
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Propanoic acid, 2-methyl-, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.263	40.05 ng/ul	14551700	Acenaphthene-d10	14.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-, 3-hyd...	216	C12H24O3	000077-68-9	90
2	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	83
3	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	80
4	Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	78
5	Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	53





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Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

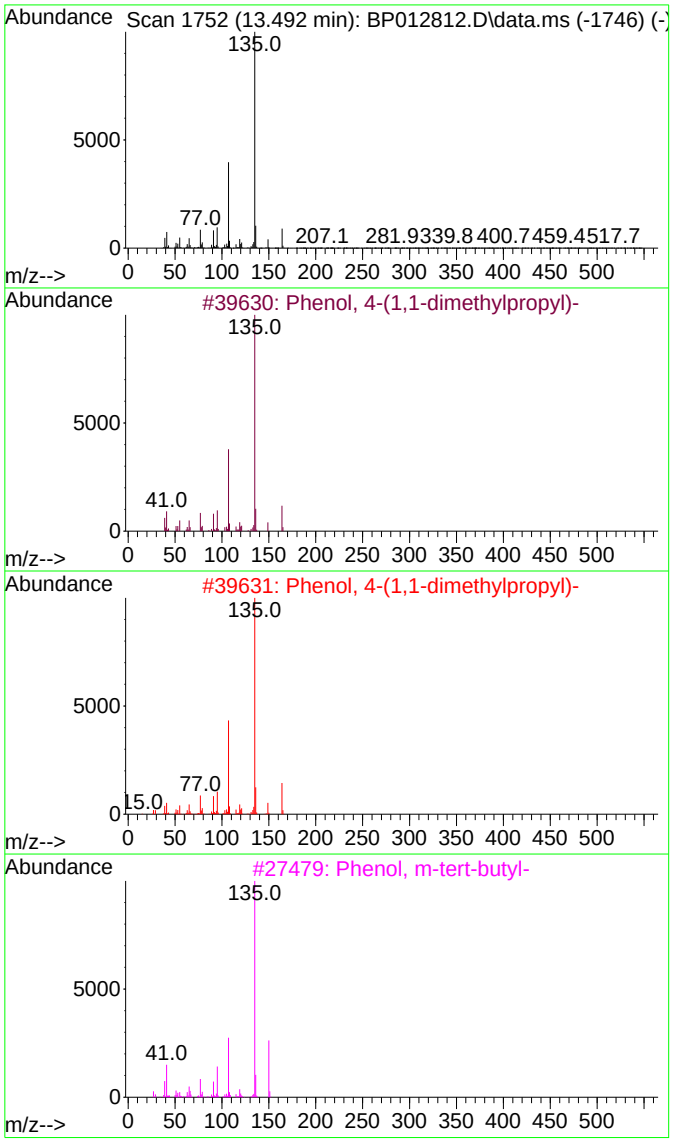
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

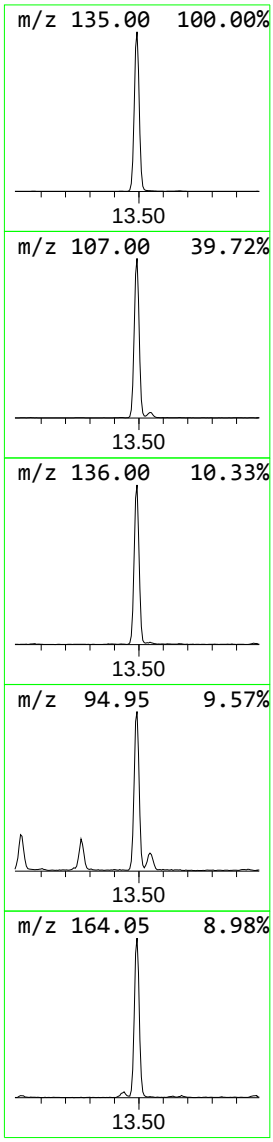
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Phenol, 4-(1,1-dimethylprop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.492	10.82 ng/ul	3930830	Acenaphthene-d10	14.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	94
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	94
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	86
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	86
5	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	81





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Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

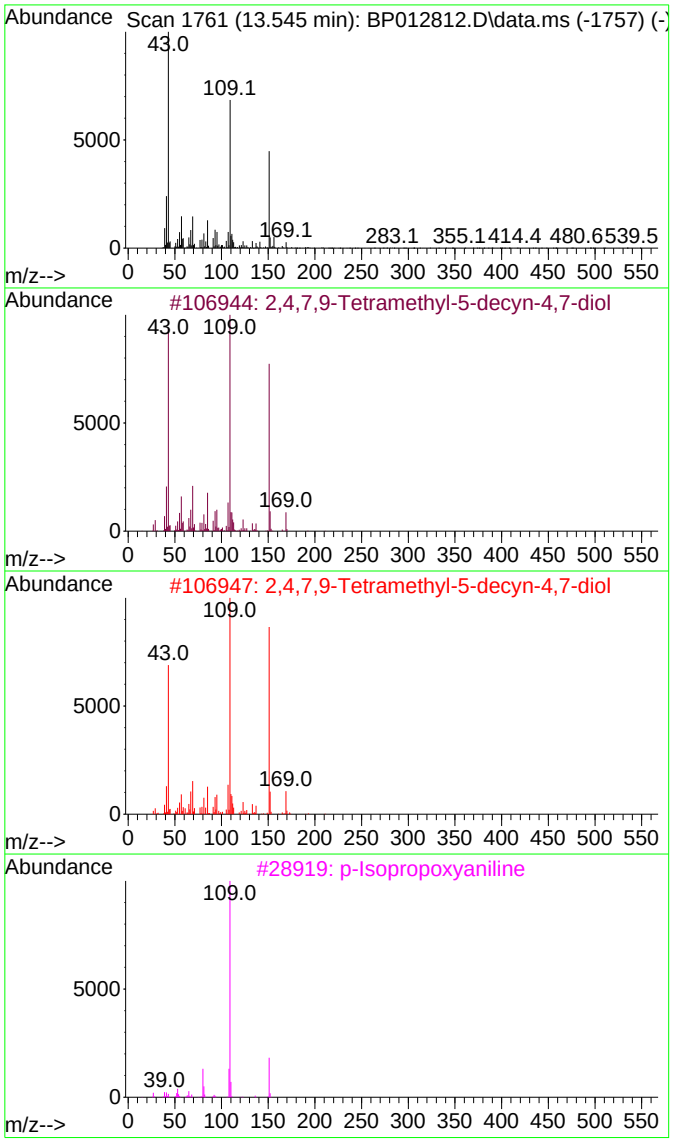
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

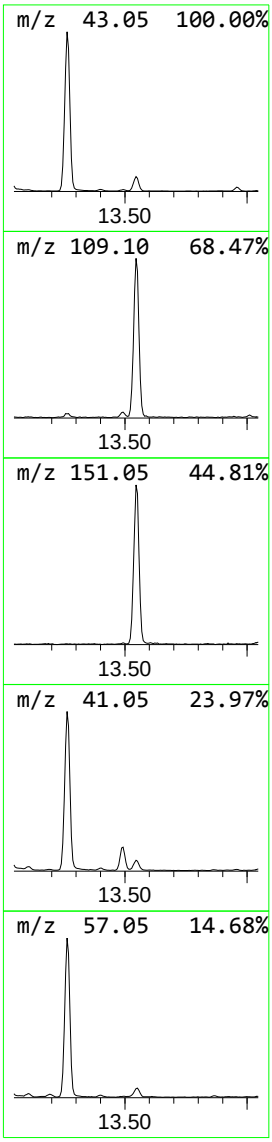
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.545	2.70 ng/ul	980397	Acenaphthene-d10	14.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91
2	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	90
3	p-Isopropoxyaniline	151	C9H13NO	007664-66-6	58
4	Acetaminophen	151	C8H9NO2	000103-90-2	53
5	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	53





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Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

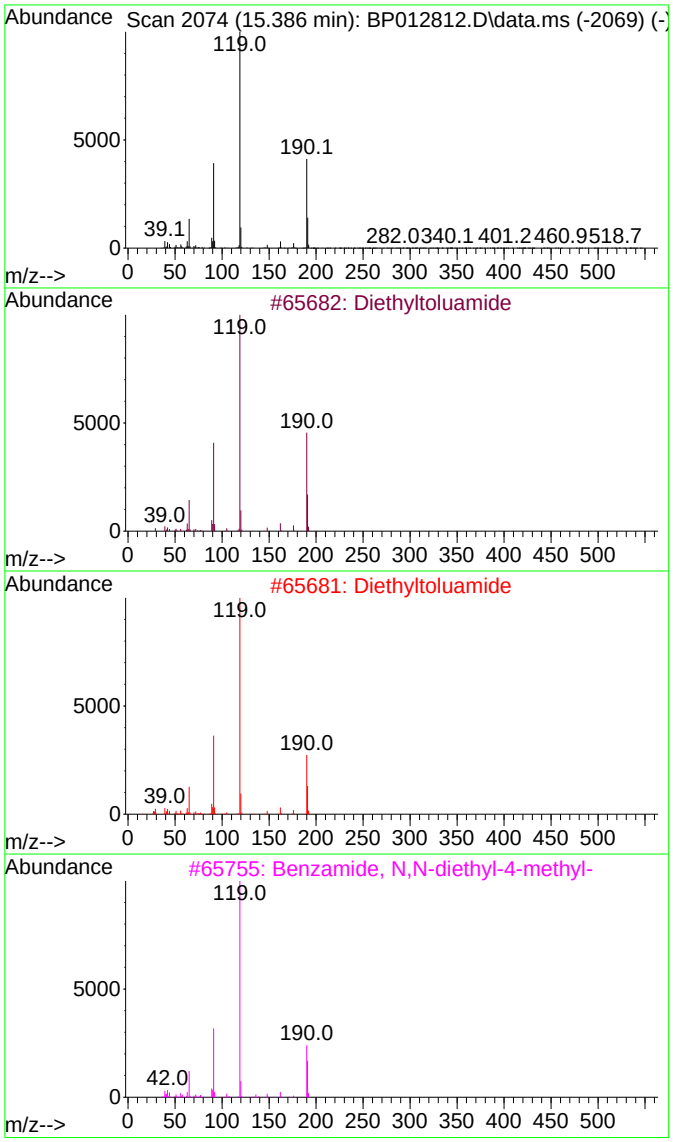
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

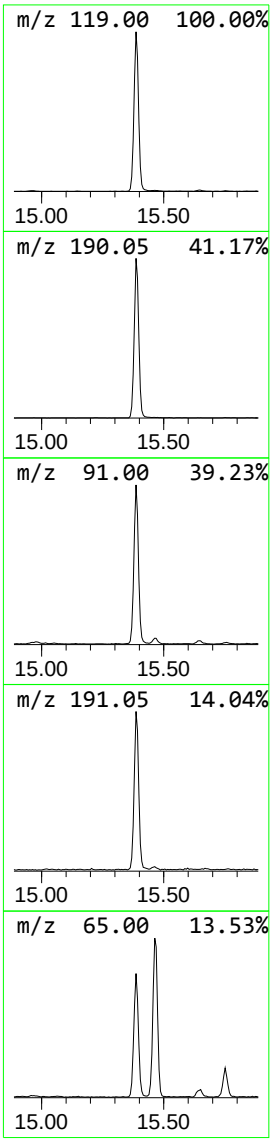
****

Peak Number 11 Diethyltoluamide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.387	6.10 ng/ul	2216980	Acenaphthene-d10	14.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	96
2	Diethyltoluamide	191	C12H17NO	000134-62-3	94
3	Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	76
4	Benzamide, 4-methyl-N-butyl-	191	C12H17NO	1000407-46-6	76
5	Benzamide, 3-methyl-N-butyl-	191	C12H17NO	1000407-41-1	76





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Acq On : 28 Nov 2022 12:27
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Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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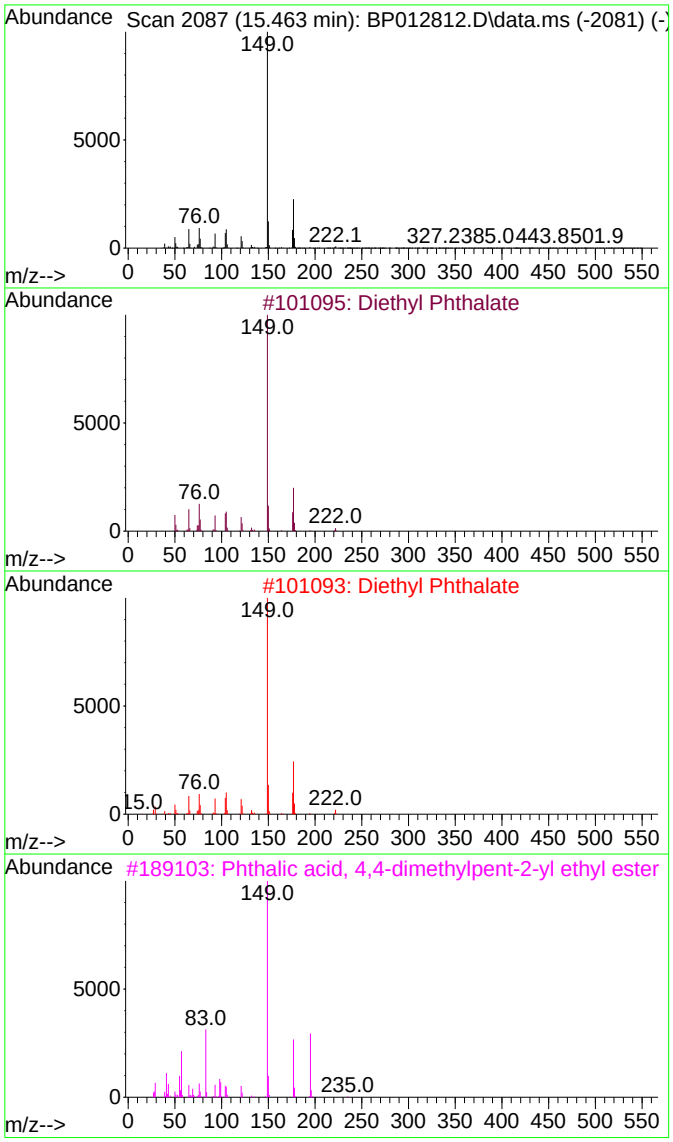
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

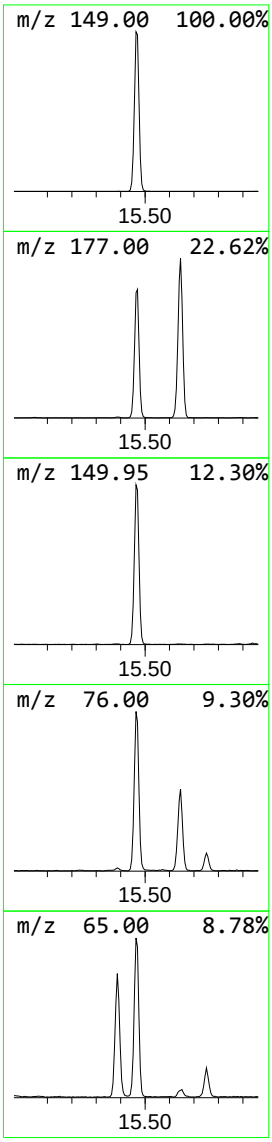
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Diethyl Phthalate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.463	10.57 ng/ul	3839310	Acenaphthene-d10	14.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyl Phthalate	222	C12H14O4	000084-66-2	93
2	Diethyl Phthalate	222	C12H14O4	000084-66-2	91
3	Phthalic acid, 4,4-dimethylpent-...	292	C17H24O4	1000371-14-4	86
4	Diethyl Phthalate	222	C12H14O4	000084-66-2	83
5	Phthalic acid, 7-bromoheptyl eth...	370	C17H23BrO4	1000415-51-6	83





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Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

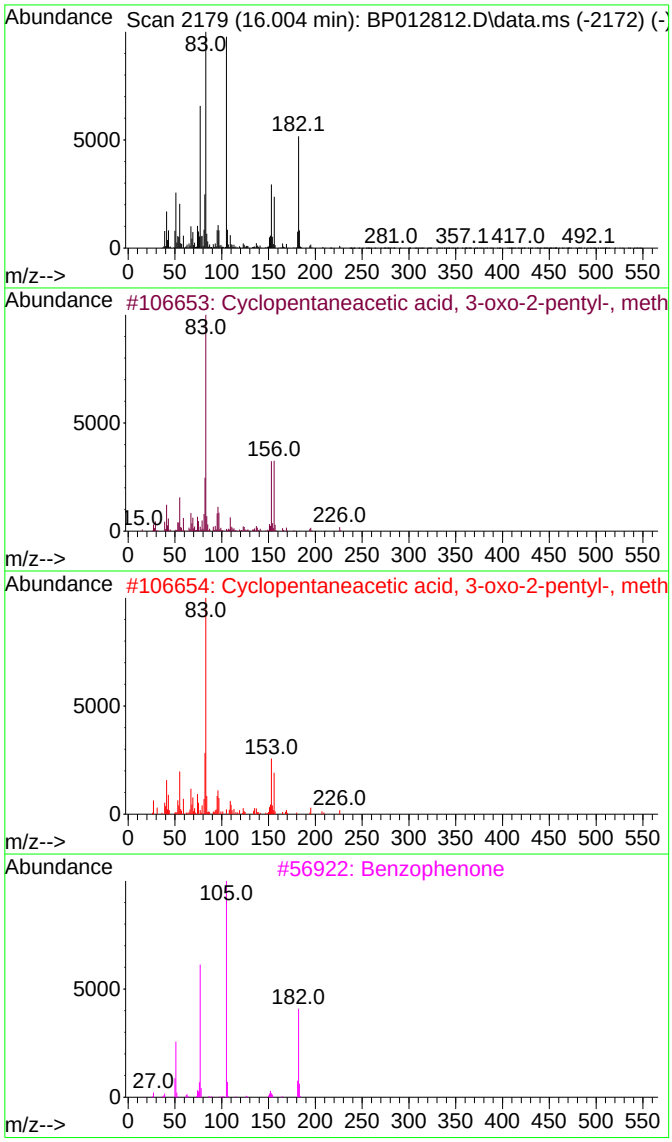
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

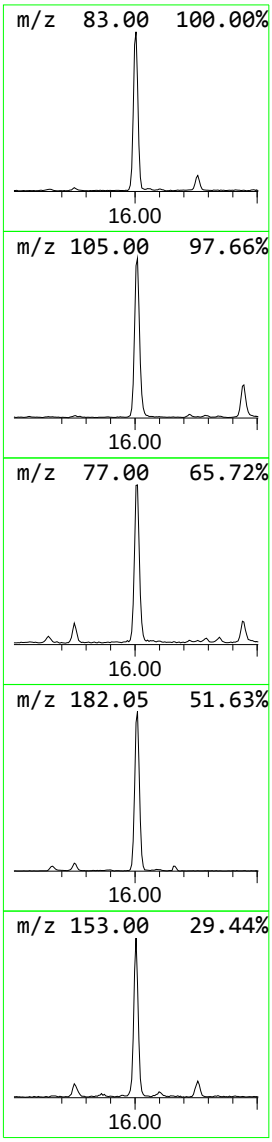
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Cyclopentaneacetic acid, 3-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.004	5.00 ng/ul	1815650	Acenaphthene-d10	14.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentaneacetic acid, 3-oxo-2...	226	C13H22O3	024851-98-7	91
2	Cyclopentaneacetic acid, 3-oxo-2...	226	C13H22O3	024851-98-7	70
3	Benzophenone	182	C13H10O	000119-61-9	46
4	Benzophenone	182	C13H10O	000119-61-9	46
5	Benzophenone	182	C13H10O	000119-61-9	46





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Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

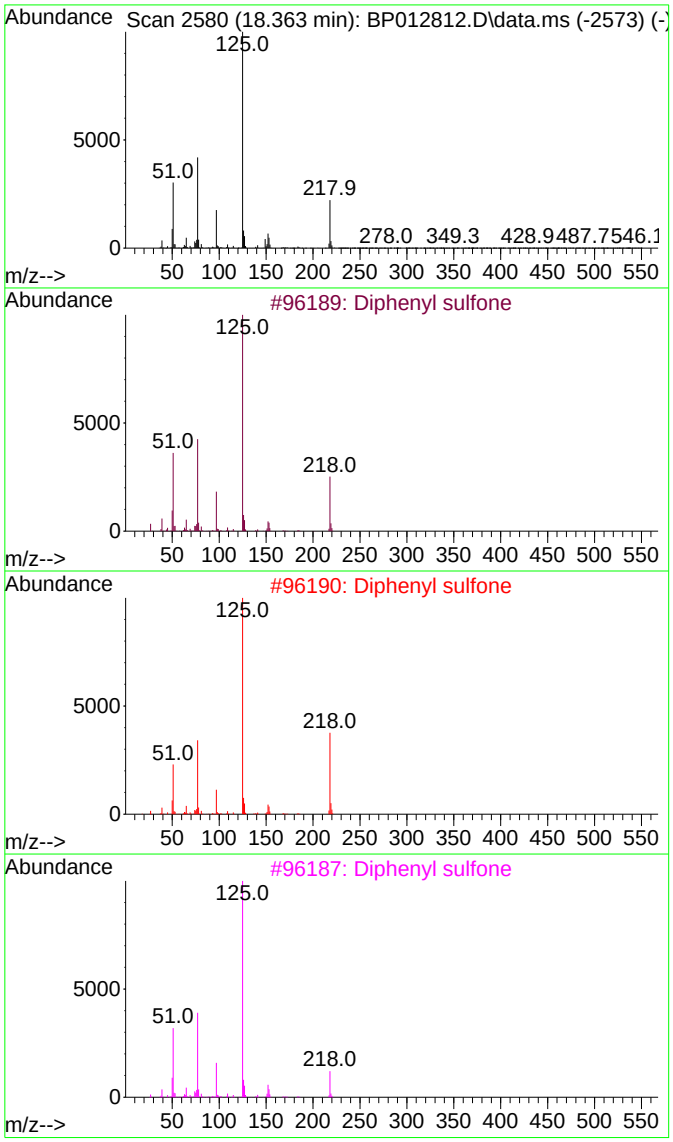
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

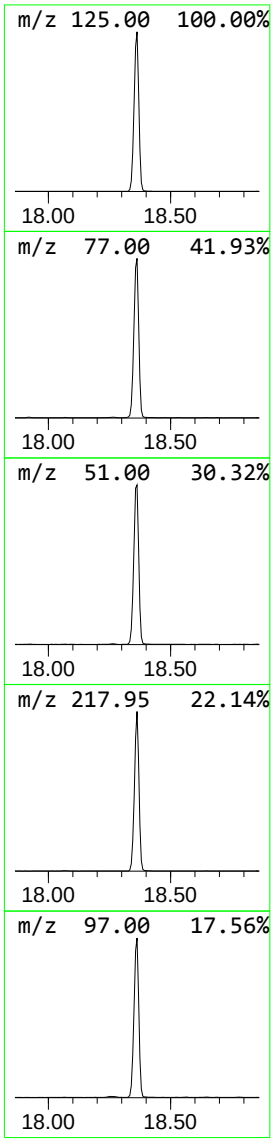
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Diphenyl sulfone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.363	30.41 ng/ul	13139100	Phenanthrene-d10	17.404

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diphenyl sulfone	218	C12H10O2S	000127-63-9	98
2	Diphenyl sulfone	218	C12H10O2S	000127-63-9	97
3	Diphenyl sulfone	218	C12H10O2S	000127-63-9	96
4	Diphenyl sulfone	218	C12H10O2S	000127-63-9	70
5	2-Propenenitrile, 3-(phenylsulfo...	193	C9H7NO2S	001424-51-7	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012812.D
Acq On : 28 Nov 2022 12:27
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.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
Ethanol, 2-butoxy-	6.075	2.2	ng/ul	338063	1	8.016	3063390	20.0
2-Propanol, 1-b...	6.652	2.4	ng/ul	369364	1	8.016	3063390	20.0
Ethanol, 2-(hex...	9.346	3.3	ng/ul	497844	1	8.016	3063390	20.0
Ethanol, 2-(2-b...	10.604	3.9	ng/ul	937623	2	10.834	4859900	20.0
Ethanol, 2-phen...	11.187	2.7	ng/ul	664286	2	10.834	4859900	20.0
Benzothiazole	11.481	2.5	ng/ul	595833	2	10.834	4859900	20.0
2,2,4-Trimethyl...	13.016	28.5	ng/ul	10348000	3	14.651	7266650	20.0
Propanoic acid,...	13.263	40.0	ng/ul	14551700	3	14.651	7266650	20.0
Phenol, 4-(1,1-...	13.492	10.8	ng/ul	3930830	3	14.651	7266650	20.0
2,4,7,9-Tetrame...	13.545	2.7	ng/ul	980397	3	14.651	7266650	20.0
Diethyltoluamide	15.387	6.1	ng/ul	2216980	3	14.651	7266650	20.0
Diethyl Phthalate	15.463	10.6	ng/ul	3839310	3	14.651	7266650	20.0
Cyclopentaneace...	16.004	5.0	ng/ul	1815650	3	14.651	7266650	20.0
Diphenyl sulfone	18.363	30.4	ng/ul	13139100	4	17.404	8642320	20.0