LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
Data File : BP012792.D
Acq On : 27 Nov 2022 05:31
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 66 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: BP012792.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	31	35	40	rBV	103112	138857	0.78%	0.061%
2	3.822	105	108	123	rBV	480460	783917	4.38%	0.343%
3	4.269	177	184	193	rBV2	80855	162866	0.91%	0.071%
4	6.075	483	491	501	rBV	217950	358520	2.00%	0.157%
5	6.652	583	589	595	rVV	275595	408022	2.28%	0.179%
6	7.163	669	676	687	rVB2	515382	1081973	6.05%	0.473%
7	7.340	699	706	718	rVB	2030499	3156657	17.64%	1.381%
8	7.540	732	740	747	rBV	1813410	3054458	17.07%	1.337%
9	7.605	747	751	760	rVB	68817	135644	0.76%	0.059%
10	7.793	777	783	791	rVV	137949	246228	1.38%	0.108%
11	8.016	814	821	827	rBV	1998855	3141790	17.55%	1.375%
12	8.075	827	831	838	rVV	119790	177703	0.99%	0.078%
13	8.716	933	940	950	rVV	1581609	2568702	14.35%	1.124%
14	8.810	950	956	957	rVV	75103	112971	0.63%	0.049%
15	8.840	957	961	966	rVB	173307	279193	1.56%	0.122%
16	9.122	1003	1009	1013	rBV2	69378	123504	0.69%	0.054%
17	9.181	1013	1019	1027	rVV	2203951	3599479	20.11%	1.575%
18	9.269	1027	1034	1041	rVV4	110788	224199	1.25%	0.098%
19	9.346	1041	1047	1056	rVV	356205	554050	3.10%	0.242%
20	9.904	1133	1142	1154	rBV	1556262	2682333	14.99%	1.174%
21	10.075	1157	1171	1175	rVV2	133420	470761	2.63%	0.206%
22	10.122	1175	1179	1186	rVB2	74592	141272	0.79%	0.062%
23	10.310	1204	1211	1220	rBV	196632	392462	2.19%	0.172%
24	10.446	1227	1234	1242	rBV	2800140	4585883	25.62%	2.007%
25	10.604	1254	1261	1270	rBV3	611928	1057905	5.91%	0.463%
26	10.693	1270	1276	1281	rVV2	158814	287675	1.61%	0.126%
27	10.834	1291	1300	1307	rBV	3039071	5084491	28.41%	2.225%
28	10.963	1314	1322	1333	rBV	2510009	4245457	23.72%	1.858%
29	11.187	1354	1360	1374	rVB	363636	729781	4.08%	0.319%
30	11.346	1382	1387	1395	rBV2	244418	521375	2.91%	0.228%
31	11.428	1395	1401	1405	rVV	142529	214600	1.20%	0.094%
32	11.481	1405	1410	1419	rVB	373551	650513	3.63%	0.285%
33	11.569	1419	1425	1430	rBV	118126	217087	1.21%	0.095%
34	11.628	1430	1435	1442	rVB6	57154	127147	0.71%	0.056%
35	11.722	1444	1451	1455	rBV2	78122	136816	0.76%	0.060%
36	12.081	1504	1512	1516	rBV2	205025	369368	2.06%	0.162%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
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Acq On : 27 Nov 2022 05:31
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 66 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	12.287	1542	1547	1550	rBV	75749	133844	0.75%	0.059%
38	12.334	1550	1555	1560	rVB4	128457	253379	1.42%	0.111%
39	12.410	1565	1568	1573	rVB2	61343	99884	0.56%	0.044%
40	12.487	1576	1581	1586	rBV2	65301	104380	0.58%	0.046%
41	12.557	1587	1593	1600	rVB3	114318	239584	1.34%	0.105%
42	12.846	1635	1642	1646	rBV2	185041	331798	1.85%	0.145%
43	12.893	1646	1650	1654	rVB	173624	235687	1.32%	0.103%
44	12.951	1654	1660	1665	rVB4	64787	122102	0.68%	0.053%
45	13.016	1665	1671	1678	rBV	7475024	11264897	62.94%	4.929%
46	13.104	1683	1686	1697	rVB	256110	372971	2.08%	0.163%
47	13.269	1706	1714	1727	rBV	10613924	15773493	88.13%	6.902%
48	13.398	1732	1736	1740	rBV	105069	134520	0.75%	0.059%
49	13.493	1746	1752	1757	rBV	3155167	4403838	24.60%	1.927%
50	13.545	1757	1761	1768	rVV	751839	1172287	6.55%	0.513%
51	13.963	1827	1832	1836	rVB	395390	549913	3.07%	0.241%
52	14.057	1842	1848	1854	rBV	5933012	8412559	47.00%	3.681%
53	14.181	1866	1869	1873	rVB2	101032	133937	0.75%	0.059%
54	14.298	1883	1889	1891	rBV3	158551	268697	1.50%	0.118%
55	14.345	1891	1897	1903	rVV	6657735	9897798	55.30%	4.331%
56	14.398	1903	1906	1917	rVB2	195201	348334	1.95%	0.152%
57	14.651	1936	1949	1956	rBV	5188841	7657997	42.79%	3.351%
58	14.834	1971	1980	1988	rVV	432252	700070	3.91%	0.306%
59	14.922	1990	1995	1999	rVV	80985	143411	0.80%	0.063%
60	14.981	1999	2005	2009	rVV	69266	137662	0.77%	0.060%
61	15.057	2014	2018	2027	rVB7	58639	118284	0.66%	0.052%
62	15.239	2045	2049	2054	rVB	164137	235289	1.31%	0.103%
63	15.345	2062	2067	2069	rBV	73206	105897	0.59%	0.046%
64	15.387	2069	2074	2081	rVV	1807019	2517977	14.07%	1.102%
65	15.469	2081	2088	2094	rVB	3046474	4247807	23.73%	1.859%
66	15.645	2112	2118	2130	rVB	9207727	12873322	71.92%	5.633%
67	15.751	2130	2136	2147	rBV	1820663	2510138	14.02%	1.098%
68	16.004	2173	2179	2187	rVV2	1423174	2062541	11.52%	0.903%
69	16.098	2192	2195	2200	rVB6	69575	100991	0.56%	0.044%
70	16.163	2202	2206	2212	rBV4	69815	101480	0.57%	0.044%
71	16.257	2218	2222	2225	rBV	114147	144488	0.81%	0.063%
72	16.286	2225	2227	2232	rVB2	108171	132108	0.74%	0.058%
73	16.345	2232	2237	2242	rVB	185967	259427	1.45%	0.114%
74	16.439	2250	2253	2257	rVB	110052	131159	0.73%	0.057%
75	16.651	2284	2289	2295	rBV	474125	658388	3.68%	0.288%
76	16.763	2301	2308	2315	rBV3	248463	572475	3.20%	0.251%

LSC Area Percent Report

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Acq On : 27 Nov 2022 05:31
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 66 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	rBV3	70159	120239	0.67%	0.053%
78	17.157	2372	2375	2379	rVV	283843	331739	1.85%	0.145%
79	17.198	2379	2382	2388	rVB2	133193	195901	1.09%	0.086%
80	17.269	2388	2394	2401	rBV2	107938	176344	0.99%	0.077%
81	17.357	2404	2409	2412	rBV4	71653	118624	0.66%	0.052%
82	17.410	2412	2418	2426	rVB	6088799	9007446	50.32%	3.942%
83	17.504	2428	2434	2445	rBV	9894642	14659660	81.90%	6.415%
84	17.686	2462	2465	2474	rVV	222117	341773	1.91%	0.150%
85	17.763	2474	2478	2484	rVB2	65135	112740	0.63%	0.049%
86	17.969	2510	2513	2519	rBV	84989	103888	0.58%	0.045%
87	18.069	2525	2530	2534	rVV2	245387	325765	1.82%	0.143%
88	18.269	2554	2564	2569	rBV4	188461	473253	2.64%	0.207%
89	18.357	2573	2579	2595	rVB	9504519	14011751	78.28%	6.131%
90	18.645	2624	2628	2632	rBV5	115569	155332	0.87%	0.068%
91	18.780	2647	2651	2658	rVV5	84243	157695	0.88%	0.069%
92	18.845	2659	2662	2670	rVB	104183	136871	0.76%	0.060%
93	19.310	2737	2741	2747	rVB2	132802	175528	0.98%	0.077%
94	19.375	2748	2752	2757	rVB	149731	193460	1.08%	0.085%
95	19.439	2758	2763	2769	rVB6	138922	204555	1.14%	0.090%
96	19.504	2771	2774	2778	rBV2	97274	123185	0.69%	0.054%
97	19.733	2807	2813	2825	rVV	13266855	17898620	100.00%	7.832%
98	21.492	3106	3112	3124	rBV	7757560	10177660	56.86%	4.454%
99	23.845	3504	3512	3520	rBV2	8369480	17351568	96.94%	7.593%
100	24.010	3532	3540	3553	rVB2	4603671	10080613	56.32%	4.411%

Sum of corrected areas: 228522682

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
Data File : BP012792.D
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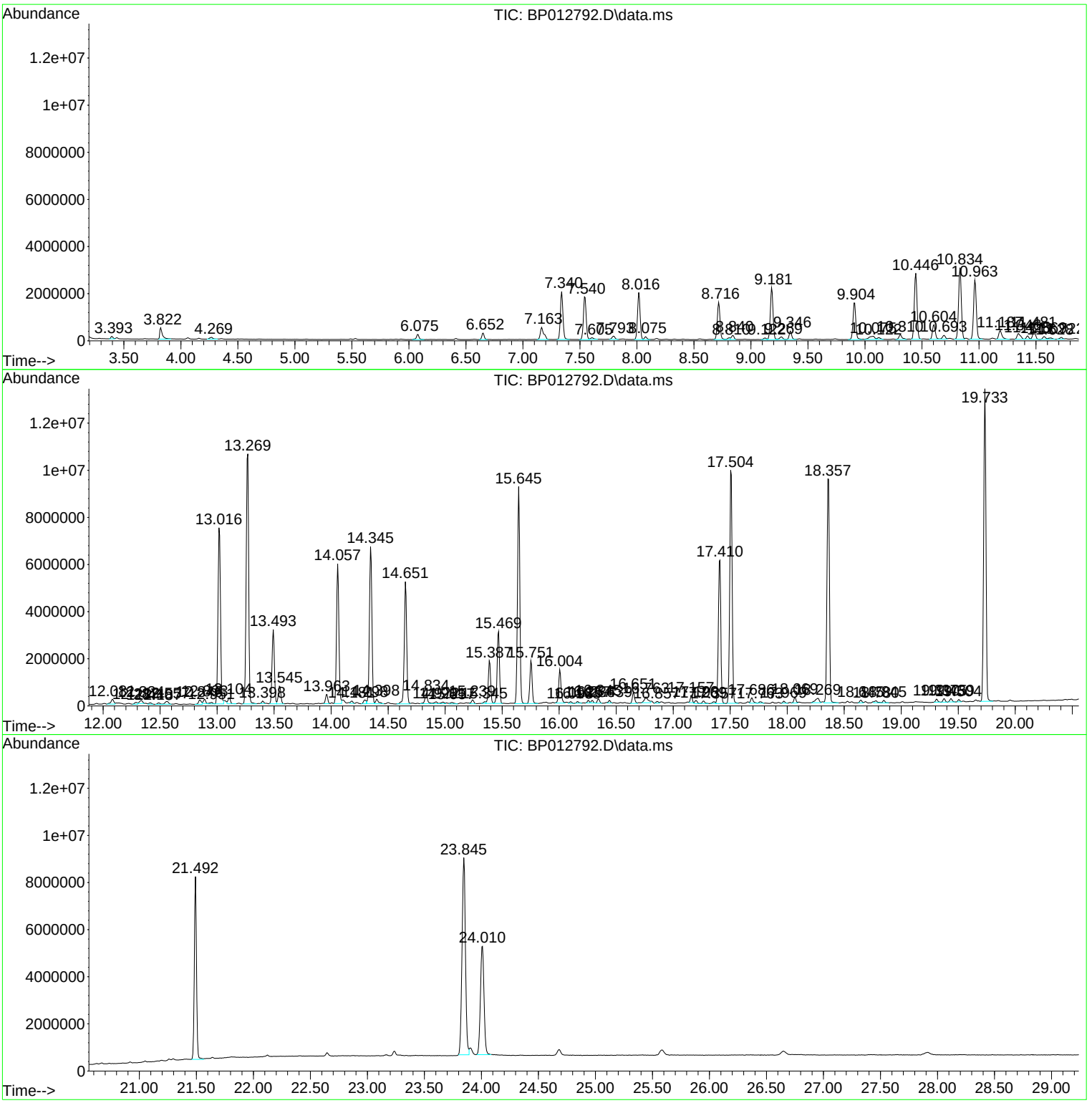
Sample : N5710-17
Misc :
ALS Vial : 66 Sample Multiplier: 1

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\BP112522\
Data File : BP012792.D
Acq On : 27 Nov 2022 05:31
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 66 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

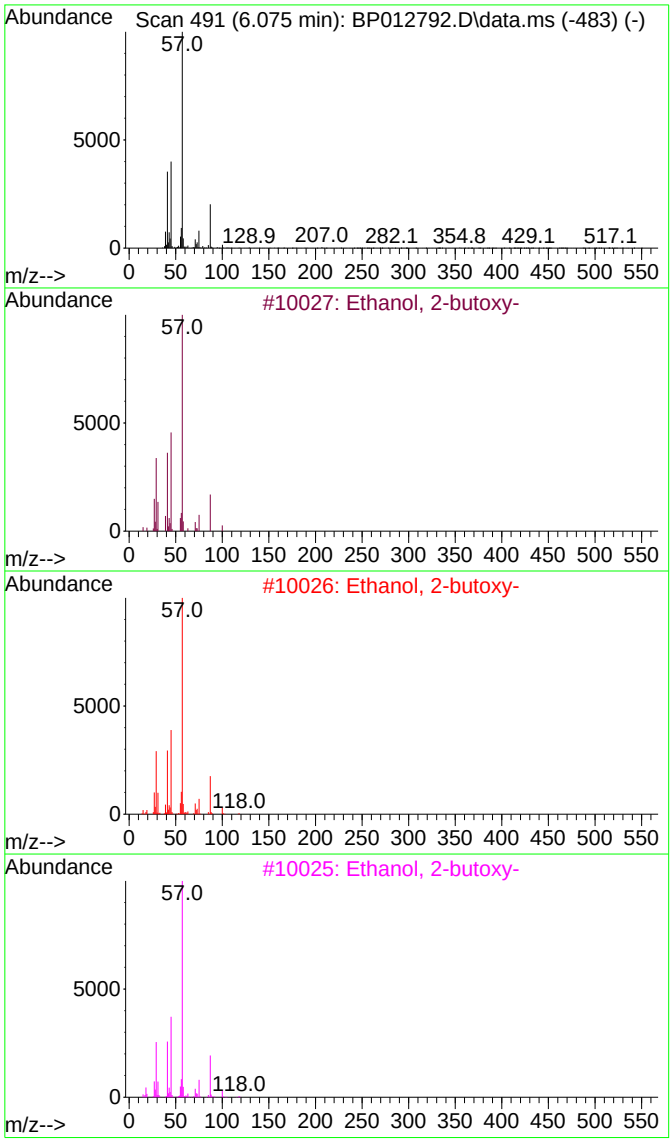
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

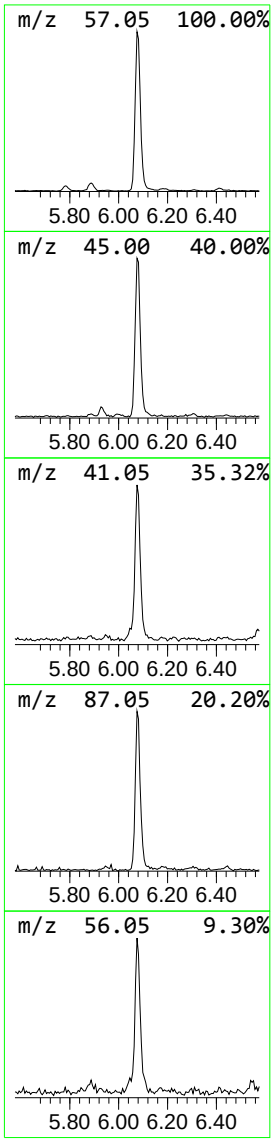
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.075	2.28 ng/ul	358520	1,4-Dichlorobenzene-d4	8.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	83
2	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	83
3	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	74
4	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	64
5	Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
Data File : BP012792.D
Acq On : 27 Nov 2022 05:31
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 66 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

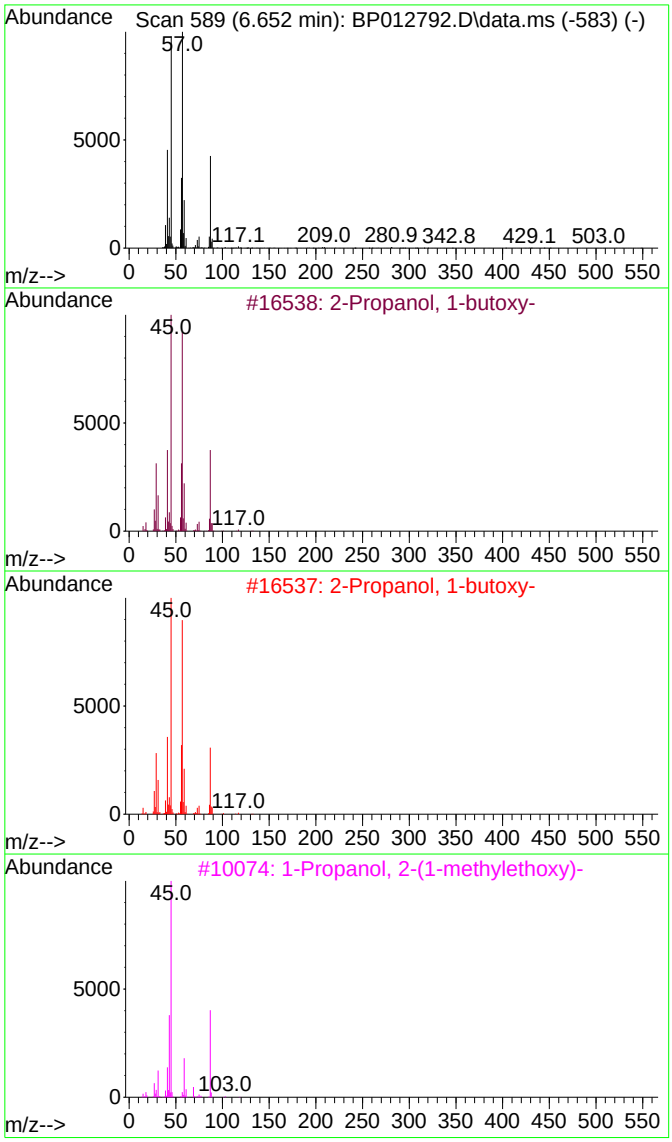
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

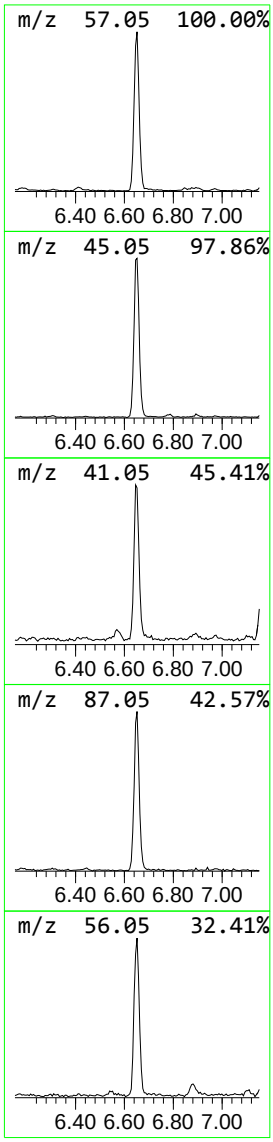
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Propanol, 1-butoxy- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.652	2.60 ng/ul	408022	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	53
4	1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	53
5	3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
Data File : BP012792.D
Acq On : 27 Nov 2022 05:31
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Sample : N5710-17
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ALS Vial : 66 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

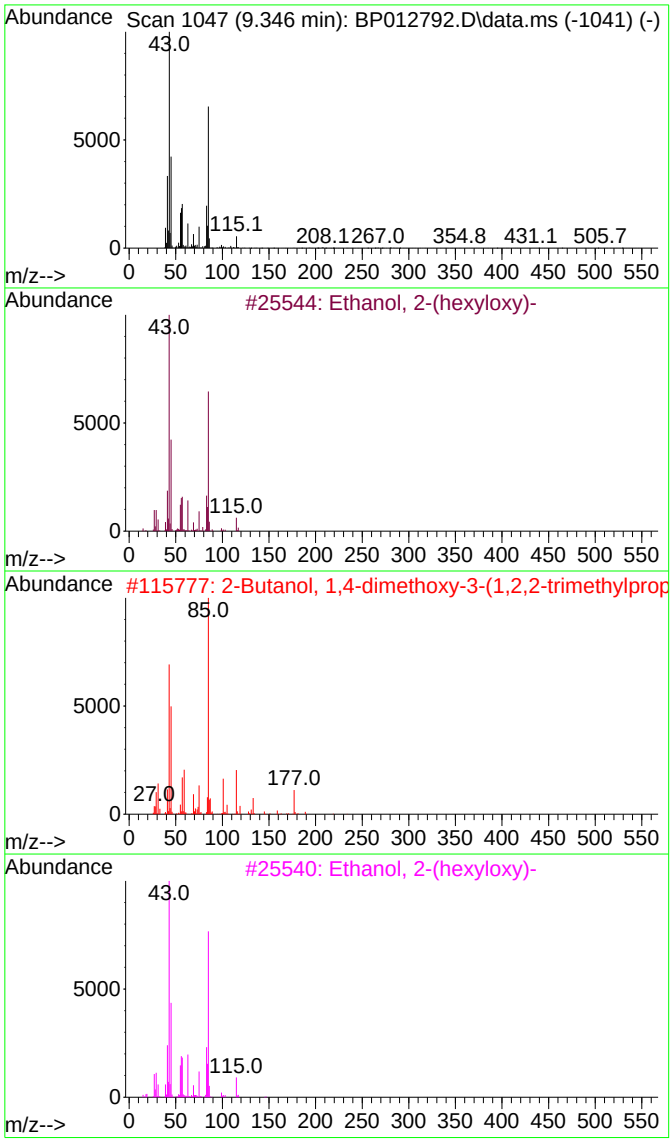
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

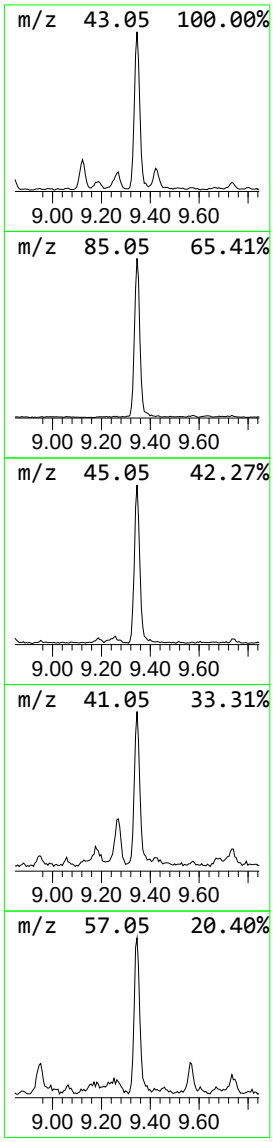
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.346	3.53 ng/ul	554050	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	2-Butanol, 1,4-dimethoxy-3-(1,2,...	234	C12H26O4	079449-92-6	50
3	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	50
4	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	42
5	Hexanal dihexyl acetal	286	C18H38O2	1000430-99-9	37





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
Data File : BP012792.D
Acq On : 27 Nov 2022 05:31
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 66 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

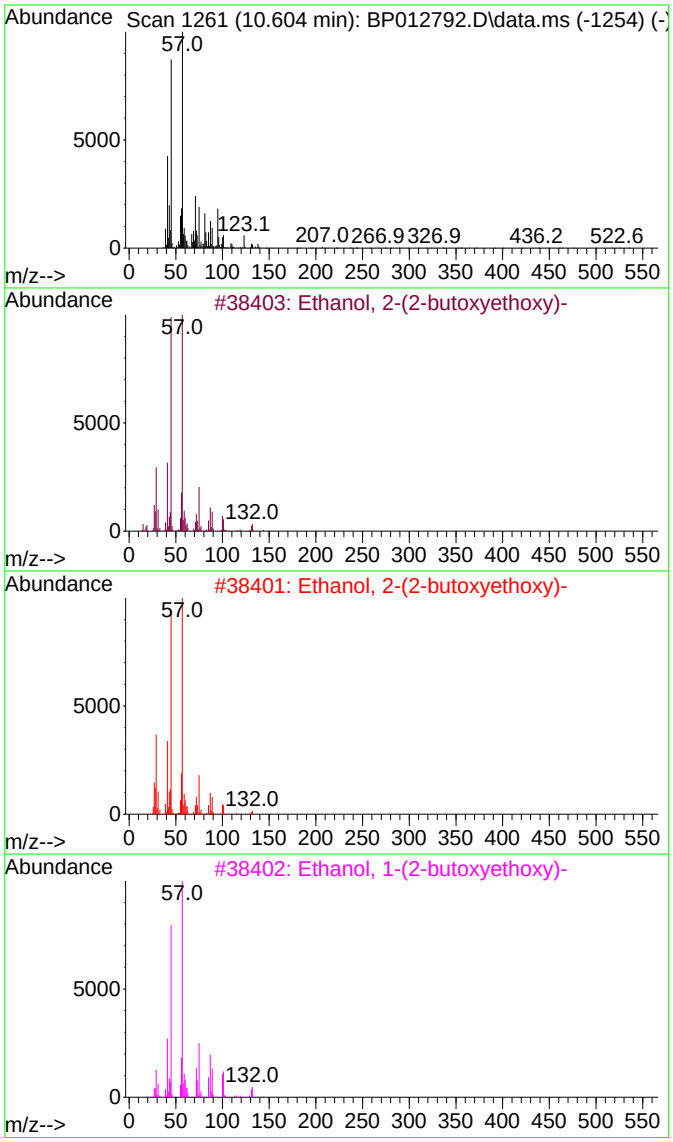
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

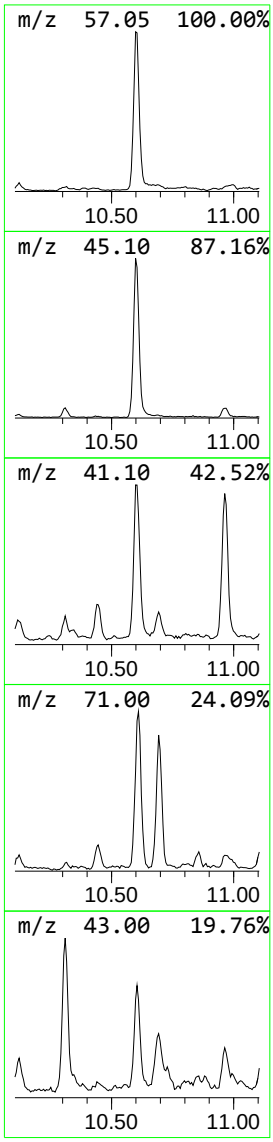
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.604	4.16 ng/ul	1057910	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
3	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	72
4	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	64
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
Data File : BP012792.D
Acq On : 27 Nov 2022 05:31
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 66 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

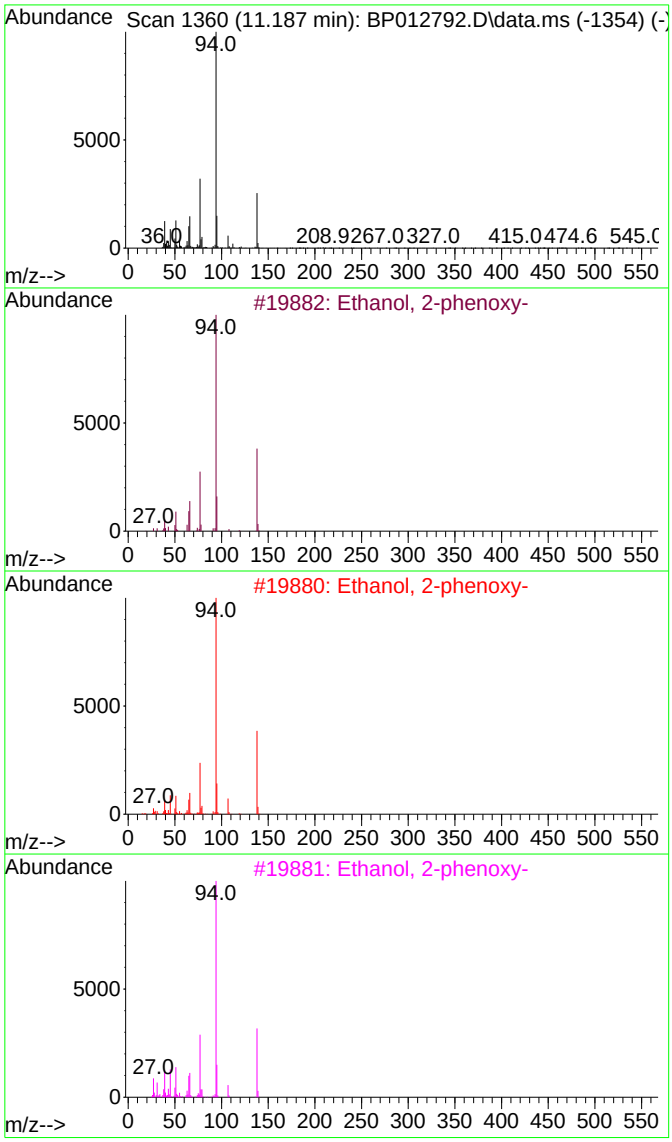
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

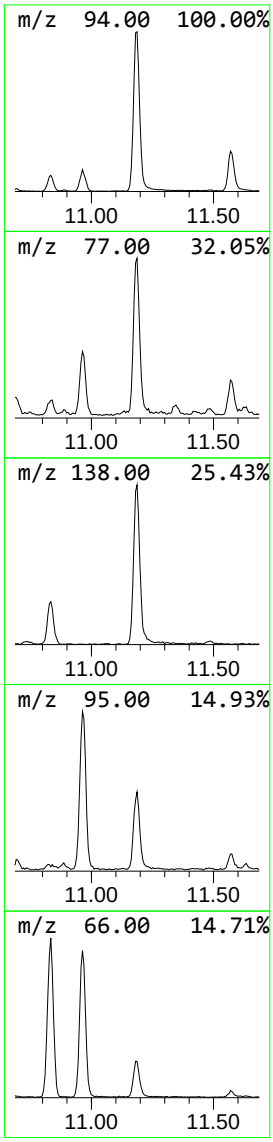
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Ethanol, 2-phenoxy- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.187	2.87 ng/ul	729781	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	94
2	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	91
3	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	91
4	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	76
5	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	64





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Acq On : 27 Nov 2022 05:31
Operator : CG/JU
Sample : N5710-17
Misc :
ALS Vial : 66 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L03

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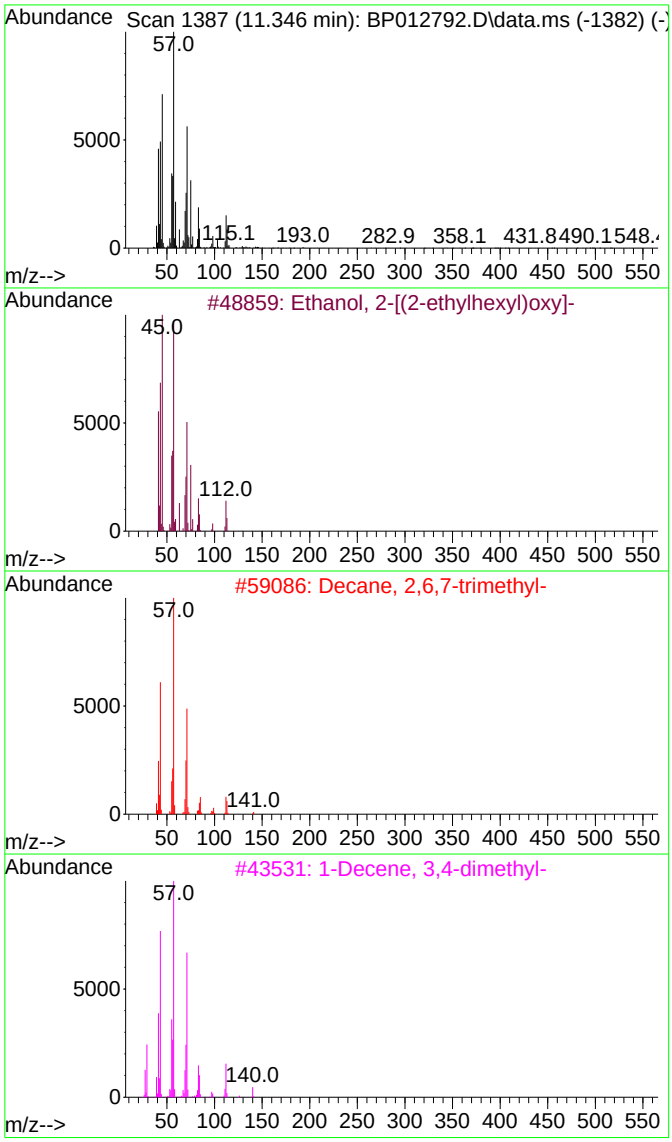
Quant Title : SVOA CALIBRATION

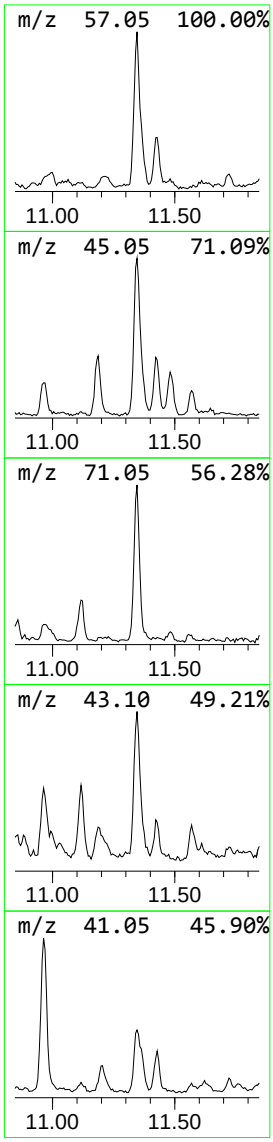
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Ethanol, 2-[(2-ethylhexyl)o... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.346	2.05 ng/ul	521375	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-[(2-ethylhexyl)oxy]-	174	C10H22O2	001559-35-9	59
2	Decane, 2,6,7-trimethyl-	184	C13H28	062108-25-2	47
3	1-Decene, 3,4-dimethyl-	168	C12H24	050871-03-9	47
4	Oxalic acid, decyl 2-ethylhexyl ...	342	C20H38O4	1000309-39-3	46
5	Carbonic acid, octyl prop-1-en-2...	214	C12H22O3	1000382-54-0	43





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Misc :
ALS Vial : 66 Sample Multiplier: 1

Instrument :
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ClientSampleId :
C0L03

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

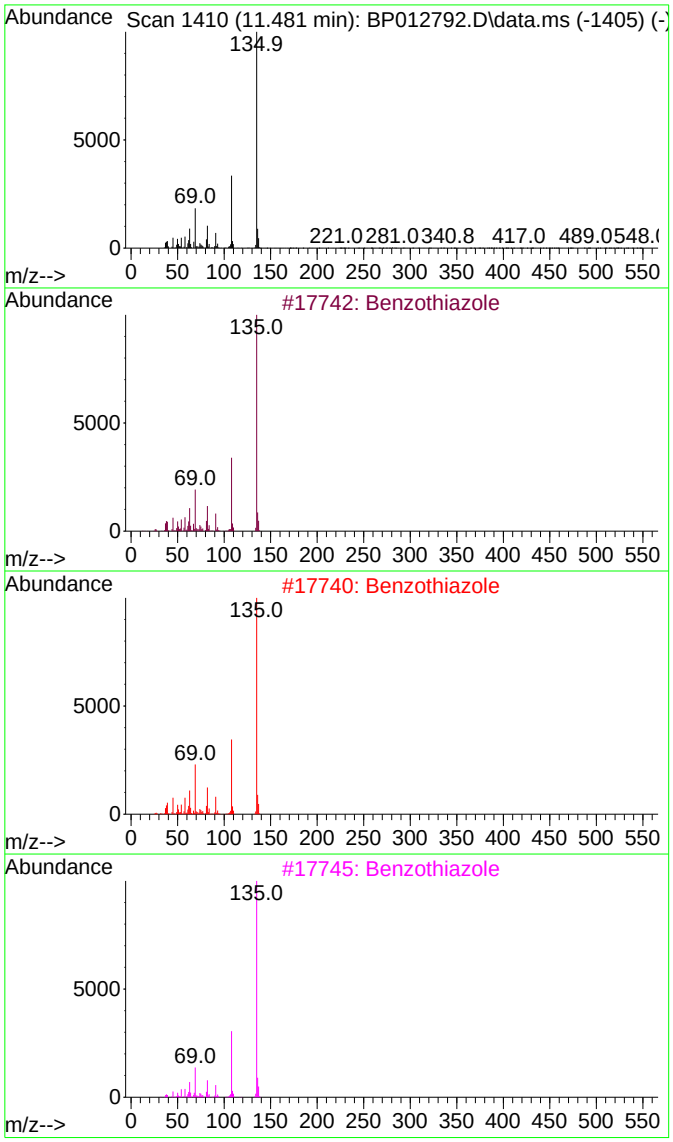
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

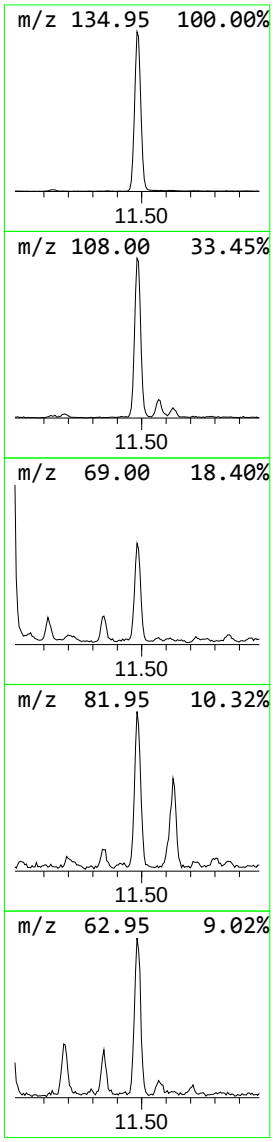
****

Peak Number 7 Benzothiazole Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.481	2.56 ng/ul	650513	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzothiazole	135	C7H5NS	000095-16-9	97
2	Benzothiazole	135	C7H5NS	000095-16-9	97
3	Benzothiazole	135	C7H5NS	000095-16-9	95
4	Benzothiazole	135	C7H5NS	000095-16-9	95
5	Benzothiazole	135	C7H5NS	000095-16-9	91





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BNA_P
ClientSampleId :
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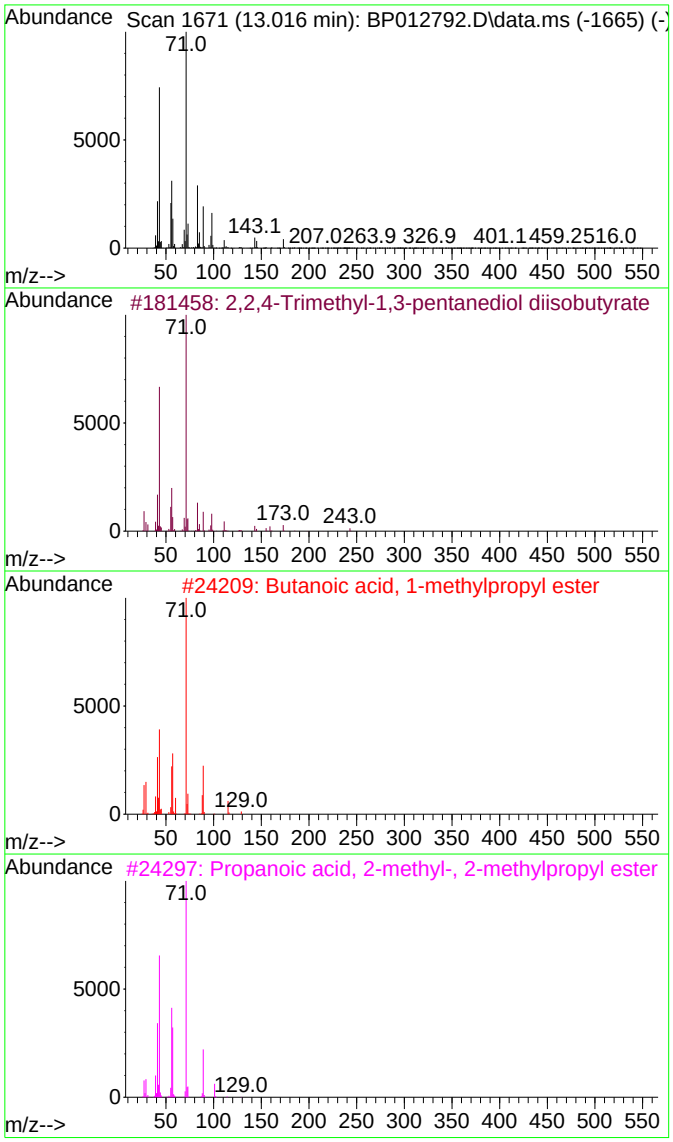
Quant Title : SVOA CALIBRATION

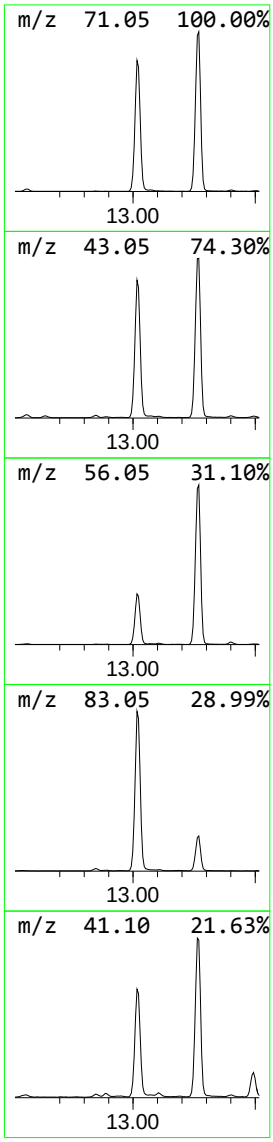
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 2,2,4-Trimethyl-1,3-pentane... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.016	29.42 ng/ul	11264900	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	56
2	Butanoic acid, 1-methylpropyl ester	144	C8H16O2	000819-97-6	42
3	Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	40
4	Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	38
5	Isophytol	296	C20H40O	000505-32-8	38





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ClientSampleId :
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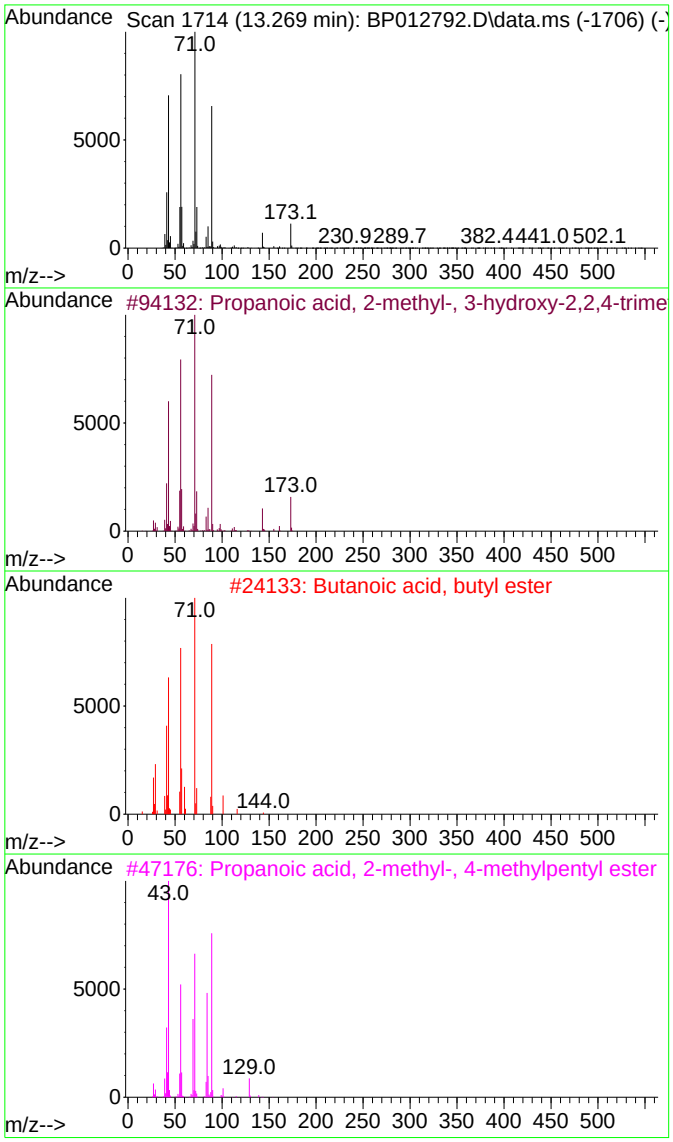
Quant Title : SVOA CALIBRATION

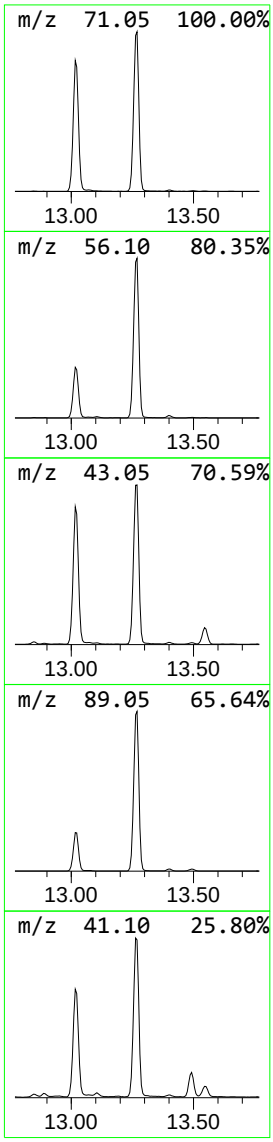
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Propanoic acid, 2-methyl-, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.269	41.19 ng/ul	15773500	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	Propanoic acid, 2-methyl-, 4-met...	172	C10H20O2	035852-44-9	56
4	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	56
5	Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	43





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Sample : N5710-17
Misc :
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ClientSampleId :
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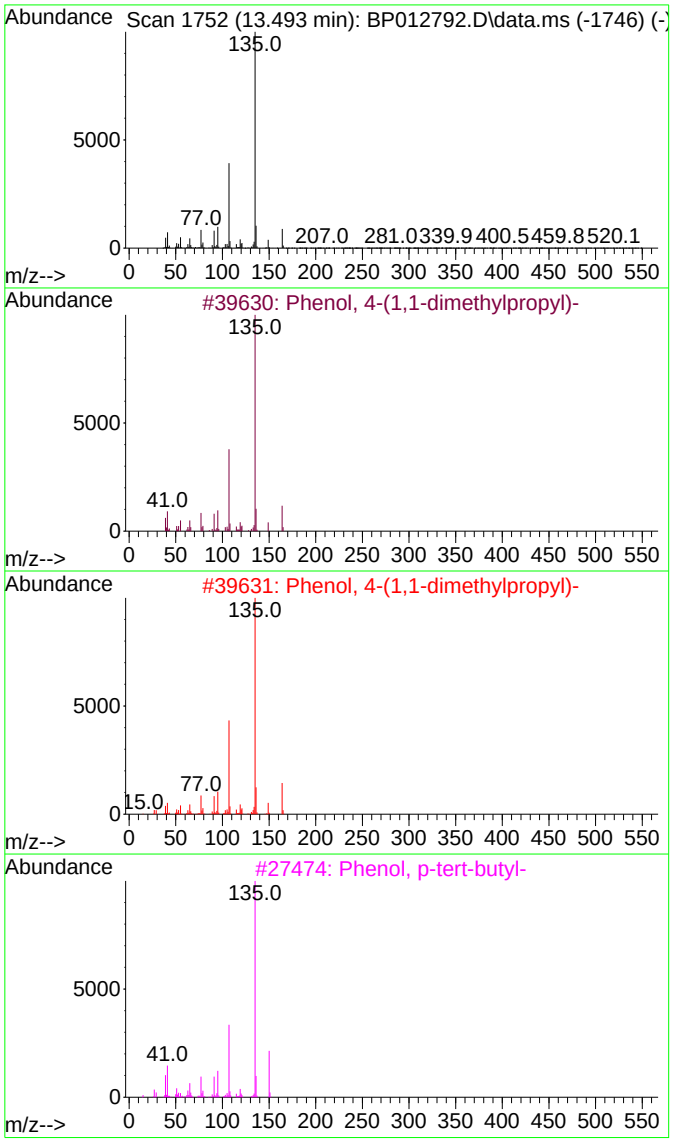
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

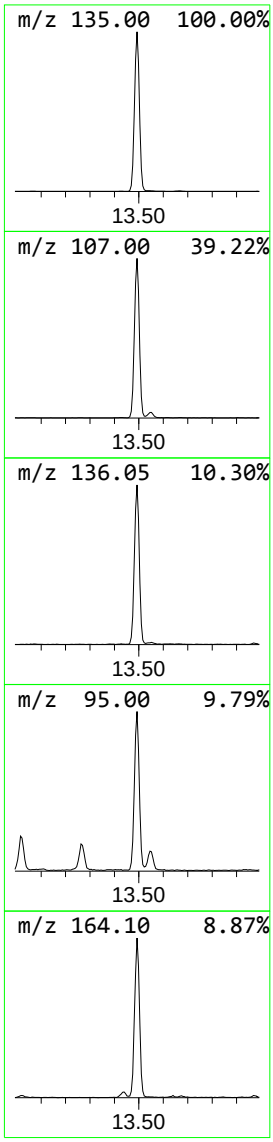
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Phenol, 4-(1,1-dimethylprop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.493	11.50 ng/ul	4403840	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	93
3	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	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 : 66 Sample Multiplier: 1

Instrument :
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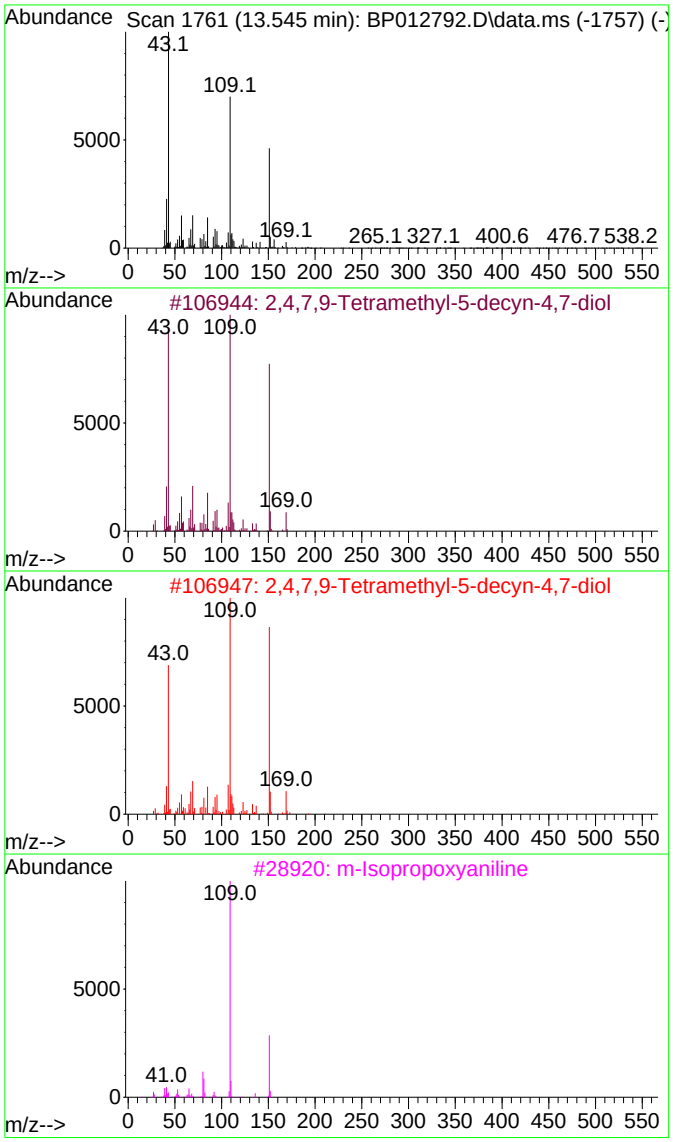
Quant Title : SVOA CALIBRATION

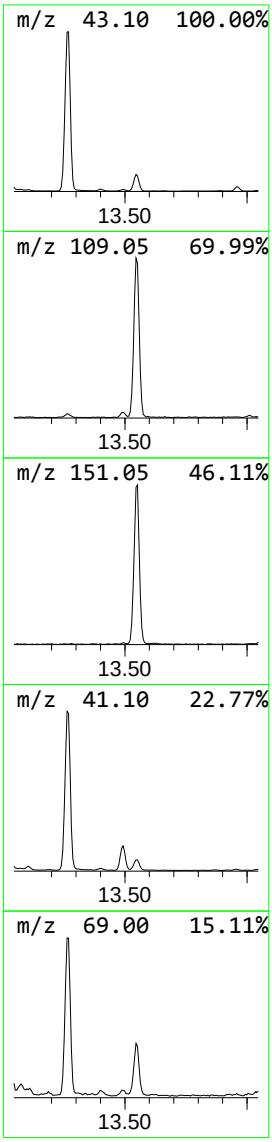
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.546	3.06 ng/ul	1172290	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	78
3	m-Isopropoxyaniline	151	C9H13NO	041406-00-2	58
4	Acetaminophen	151	C8H9NO2	000103-90-2	53
5	2-Propenal, 3-(2,2,6-trimethyl-7...	194	C12H18O2	055759-91-6	45





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Sample : N5710-17
Misc :
ALS Vial : 66 Sample Multiplier: 1

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ClientSampleId :
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Quant Title : SVOA CALIBRATION

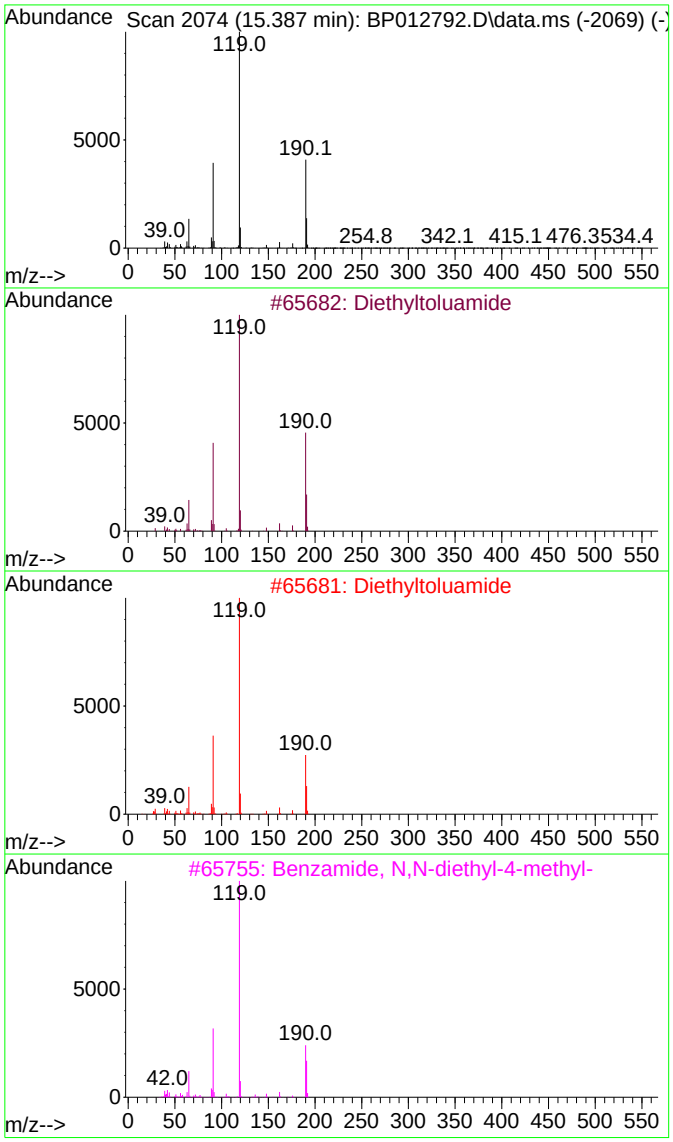
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

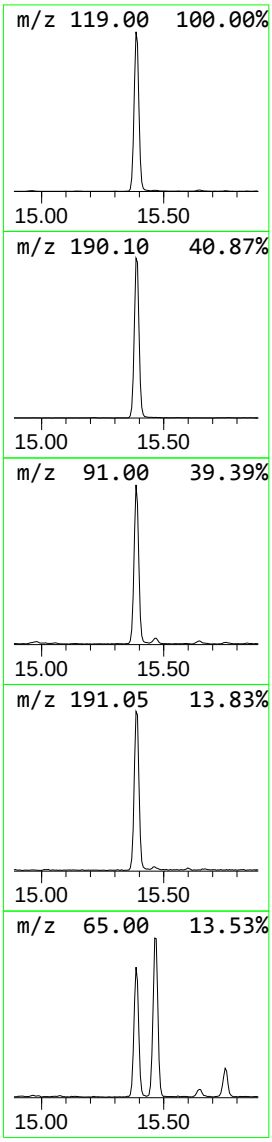
****

Peak Number 12 Diethyltoluamide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.387	6.58 ng/ul	2517980	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	87
4	Benzamide, 3-methyl-N-butyl-	191	C12H17NO	1000407-41-1	76
5	Benzamide, 3-methyl-N-methyl-N-p...	191	C12H17NO	1000421-46-2	76





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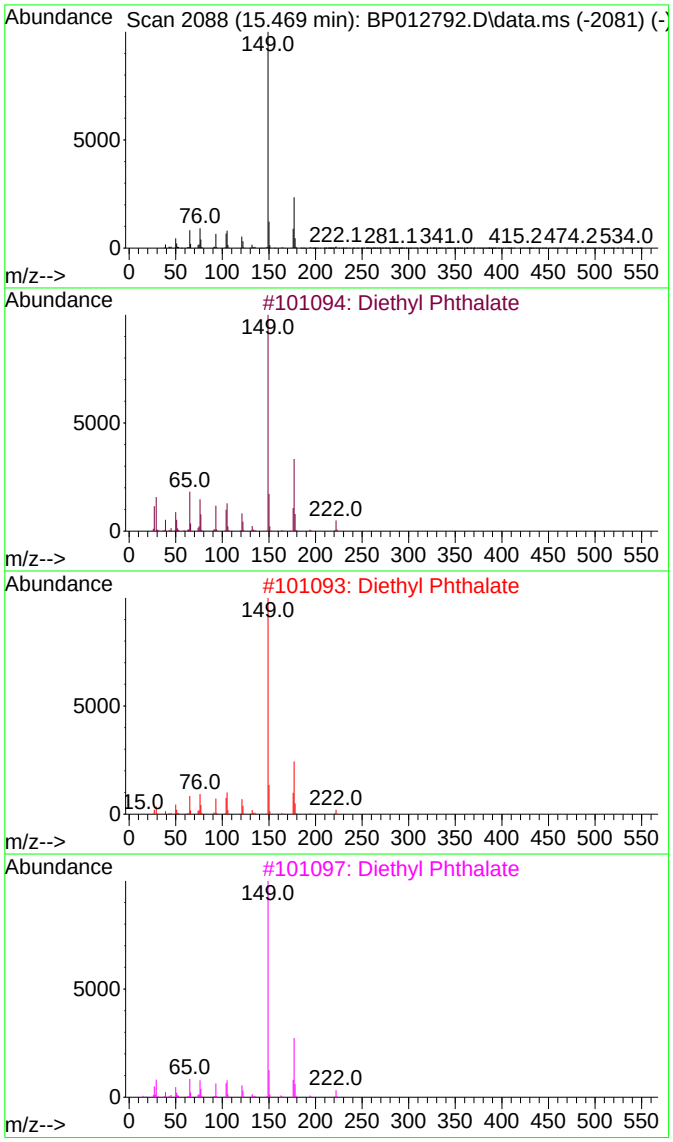
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

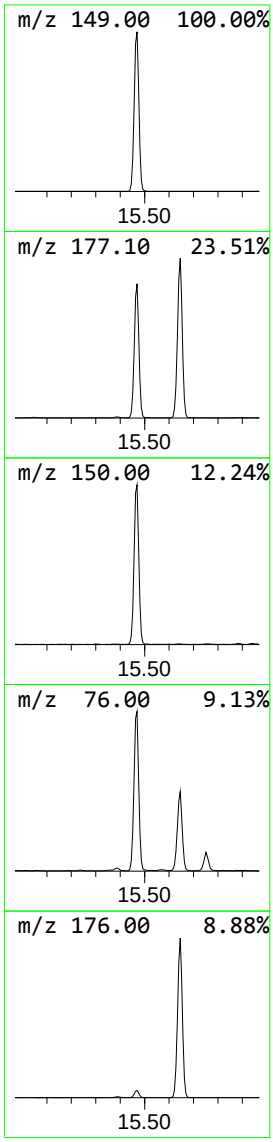
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Diethyl Phthalate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.469	11.09 ng/ul	4247810	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	Diethyl Phthalate	222	C12H14O4	000084-66-2	91
4	Diethyl Phthalate	222	C12H14O4	000084-66-2	90
5	Phthalic acid, 7-bromoheptyl eth...	370	C17H23BrO4	1000415-51-6	90





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Misc :
ALS Vial : 66 Sample Multiplier: 1

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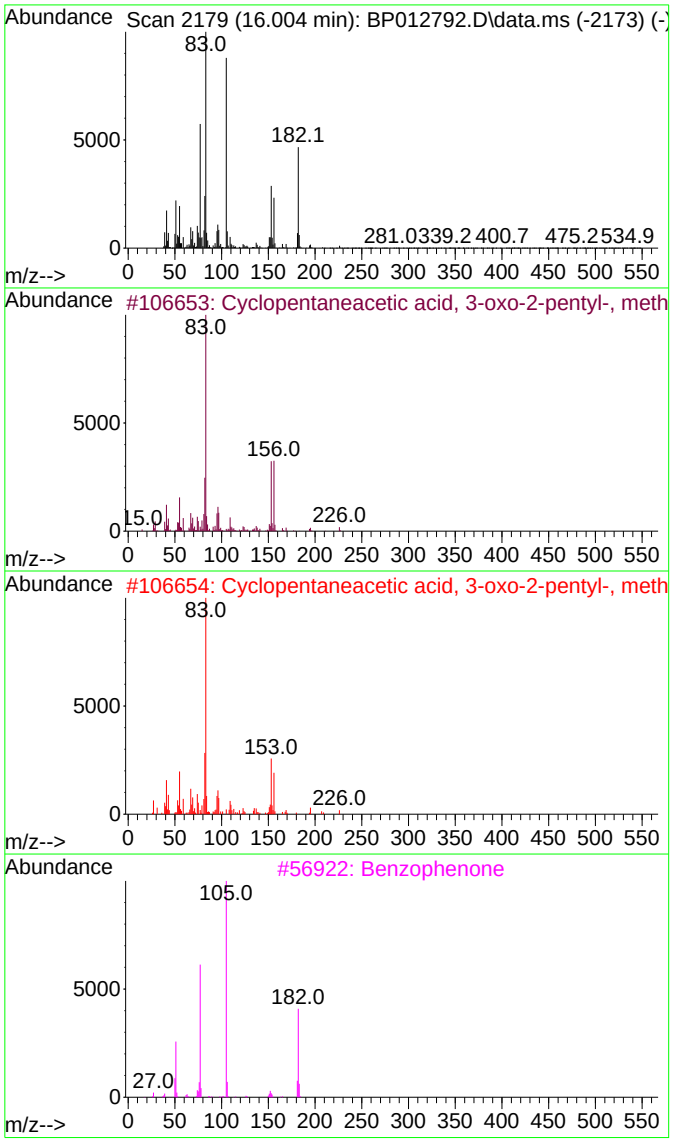
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

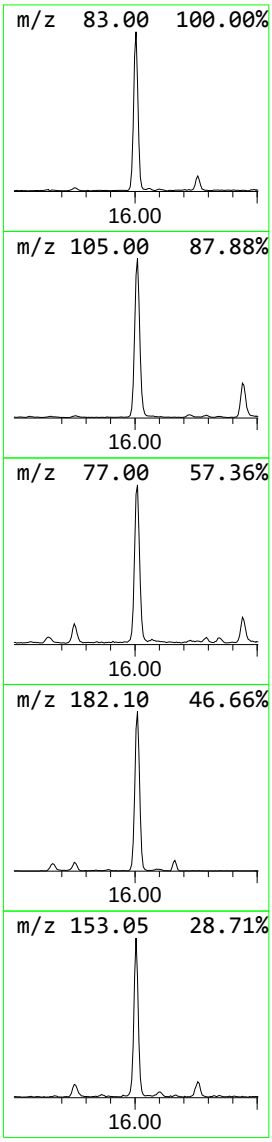
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Cyclopentaneacetic acid, 3-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.004	5.39 ng/ul	2062540	Acenaphthene-d10	14.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentaneacetic acid, 3-oxo-2...	226	C13H22O3	024851-98-7	92
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	42





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ALS Vial : 66 Sample Multiplier: 1

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BNA_P
ClientSampleId :
C0L03

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

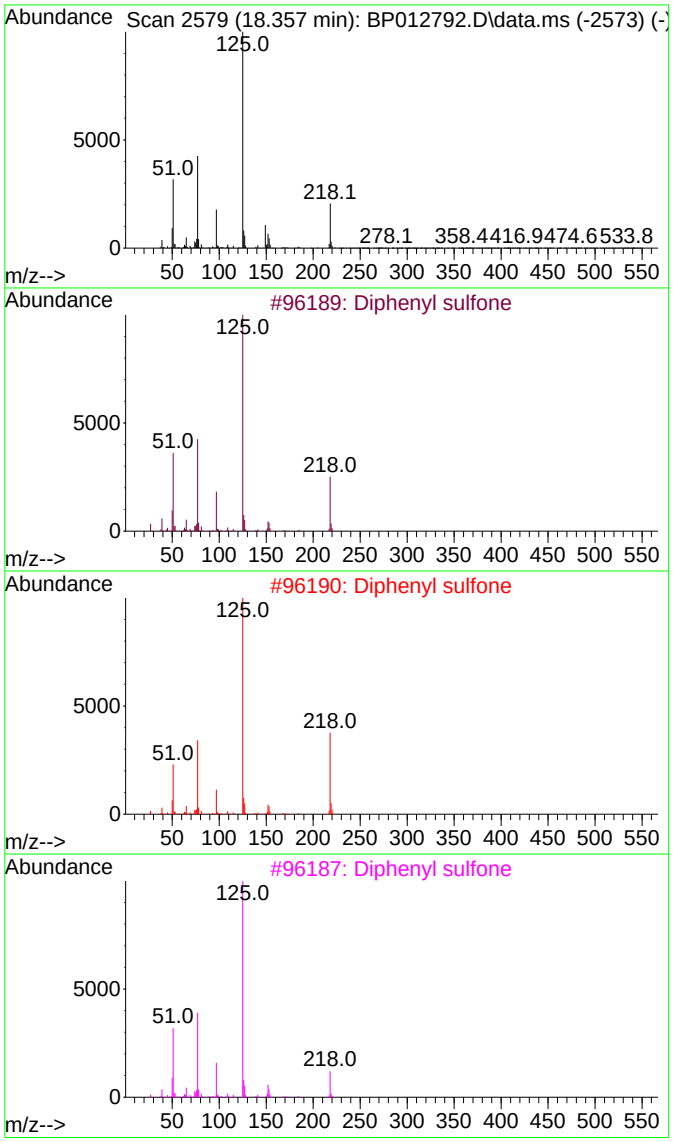
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

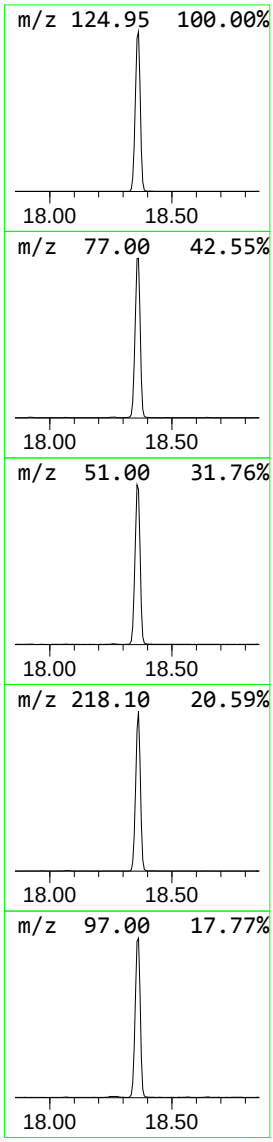
****

Peak Number 15 Diphenyl sulfone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.357	31.11 ng/ul	14011800	Phenanthrene-d10	17.410

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	Benzene, (1-propenylsulfonyl)-, ...	182	C9H10O2S	028691-72-7	50





Tentatively Identified Compound (LSC) summary

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Operator : CG/JU
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ClientSampleId :
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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	2.3	ng/ul	358520	1	8.016	3141790	20.0
2-Propanol, 1-b...	6.652	2.6	ng/ul	408022	1	8.016	3141790	20.0
Ethanol, 2-(hex...	9.346	3.5	ng/ul	554050	1	8.016	3141790	20.0
Ethanol, 2-(2-b...	10.604	4.2	ng/ul	1057910	2	10.834	5084490	20.0
Ethanol, 2-phen...	11.187	2.9	ng/ul	729781	2	10.834	5084490	20.0
Ethanol, 2-[(2-...	11.346	2.0	ng/ul	521375	2	10.834	5084490	20.0
Benzothiazole	11.481	2.6	ng/ul	650513	2	10.834	5084490	20.0
2,2,4-Trimethyl...	13.016	29.4	ng/ul	11264900	3	14.651	7658000	20.0
Propanoic acid,...	13.269	41.2	ng/ul	15773500	3	14.651	7658000	20.0
Phenol, 4-(1,1-...	13.493	11.5	ng/ul	4403840	3	14.651	7658000	20.0
2,4,7,9-Tetrame...	13.546	3.1	ng/ul	1172290	3	14.651	7658000	20.0
Diethyltoluamide	15.387	6.6	ng/ul	2517980	3	14.651	7658000	20.0
Diethyl Phthalate	15.469	11.1	ng/ul	4247810	3	14.651	7658000	20.0
Cyclopentaneace...	16.004	5.4	ng/ul	2062540	3	14.651	7658000	20.0
Diphenyl sulfone	18.357	31.1	ng/ul	14011800	4	17.410	9007450	20.0