LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012811.D
Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

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: BP012811.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	77338	108220	0.81%	0.065%
2	3.822	105	108	124	rBV	372868	621478	4.65%	0.373%
3	4.064	144	149	154	rBV	37833	59134	0.44%	0.035%
4	4.264	177	183	188	rBV2	48218	89166	0.67%	0.053%
5	6.075	487	491	498	rVB	144857	217242	1.63%	0.130%
6	6.652	583	589	596	rBV	179302	273618	2.05%	0.164%
7	7.163	669	676	688	rVB2	318005	687713	5.15%	0.413%
8	7.340	699	706	717	rVB	1397653	2209991	16.54%	1.326%
9	7.481	723	730	734	rBV2	79487	173016	1.29%	0.104%
10	7.540	734	740	748	rVV	1215654	1967976	14.73%	1.181%
11	7.605	748	751	759	rVB	35927	69083	0.52%	0.041%
12	7.793	777	783	793	rVV	88593	159827	1.20%	0.096%
13	8.016	812	821	827	rBV	1657309	2591521	19.39%	1.555%
14	8.075	827	831	837	rVB	72043	110117	0.82%	0.066%
15	8.716	932	940	951	rVV	1032969	1663751	12.45%	0.998%
16	8.804	951	955	957	rVV	48697	75406	0.56%	0.045%
17	8.840	957	961	966	rVB	116209	191133	1.43%	0.115%
18	9.122	1003	1009	1013	rBV3	48409	84567	0.63%	0.051%
19	9.181	1013	1019	1026	rVV	1493818	2388665	17.87%	1.433%
20	9.263	1027	1033	1040	rVV4	75432	158608	1.19%	0.095%
21	9.346	1041	1047	1055	rVV	223887	369274	2.76%	0.222%
22	9.904	1134	1142	1156	rBV	1043293	1721629	12.88%	1.033%
23	10.069	1157	1170	1175	rVV3	87024	286613	2.14%	0.172%
24	10.116	1175	1178	1186	rVB3	29972	65945	0.49%	0.040%
25	10.310	1204	1211	1220	rBV	129010	259415	1.94%	0.156%
26	10.446	1227	1234	1242	rBV	1949457	3139715	23.49%	1.884%
27	10.604	1255	1261	1270	rBV2	367812	662482	4.96%	0.397%
28	10.693	1270	1276	1281	rBV3	105839	182571	1.37%	0.110%
29	10.834	1290	1300	1306	rBV	2464844	4128544	30.89%	2.477%
30	10.963	1313	1322	1331	rBV	1713838	2900905	21.71%	1.740%
31	11.116	1342	1348	1354	rBV5	31121	63715	0.48%	0.038%
32	11.187	1354	1360	1373	rVB	246289	491176	3.68%	0.295%
33	11.346	1381	1387	1395	rBV2	153197	326737	2.45%	0.196%
34	11.422	1395	1400	1405	rVV	84885	133640	1.00%	0.080%
35	11.481	1405	1410	1419	rVB	257459	439249	3.29%	0.264%
36	11.569	1419	1425	1430	rBV	76257	136150	1.02%	0.082%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012811.D
Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

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	11.628	1431	1435	1442	rVB7	36522	74258	0.56%	0.045%
38	11.722	1445	1451	1454	rBV	48807	76497	0.57%	0.046%
39	12.081	1503	1512	1516	rBV2	129769	230841	1.73%	0.138%
40	12.287	1542	1547	1549	rBV	41078	64549	0.48%	0.039%
41	12.328	1550	1554	1560	rVB4	74136	151598	1.13%	0.091%
42	12.410	1565	1568	1576	rVB2	44042	73283	0.55%	0.044%
43	12.487	1576	1581	1585	rBV2	42909	70195	0.53%	0.042%
44	12.563	1586	1594	1600	rVB2	78370	161656	1.21%	0.097%
45	12.845	1635	1642	1646	rBV2	119821	229022	1.71%	0.137%
46	12.887	1646	1649	1654	rVV	120386	178289	1.33%	0.107%
47	12.951	1654	1660	1664	rVB5	41304	77749	0.58%	0.047%
48	13.016	1664	1671	1683	rBV	5180214	7855367	58.78%	4.713%
49	13.104	1683	1686	1696	rVV	185022	292069	2.19%	0.175%
50	13.263	1706	1713	1727	rVV	7850614	11195893	83.78%	6.717%
51	13.398	1731	1736	1742	rVV	77971	129801	0.97%	0.078%
52	13.492	1745	1752	1757	rBV	2182343	3074632	23.01%	1.845%
53	13.545	1757	1761	1768	rVB	484861	739139	5.53%	0.443%
54	13.957	1826	1831	1836	rVB	272110	353270	2.64%	0.212%
55	14.057	1842	1848	1854	rBV	4376258	6075631	45.46%	3.645%
56	14.181	1866	1869	1873	rVB3	69546	88135	0.66%	0.053%
57	14.228	1873	1877	1883	rVB4	40915	72070	0.54%	0.043%
58	14.298	1883	1889	1891	rBV3	114892	185022	1.38%	0.111%
59	14.345	1891	1897	1903	rVV	4848659	6803663	50.91%	4.082%
60	14.398	1903	1906	1917	rVB2	127003	238541	1.79%	0.143%
61	14.651	1937	1949	1956	rBV	4177482	6142313	45.96%	3.685%
62	14.834	1971	1980	1989	rBV	204477	373650	2.80%	0.224%
63	14.922	1989	1995	1998	rVB4	45548	71462	0.53%	0.043%
64	15.057	2014	2018	2026	rVB7	38562	81004	0.61%	0.049%
65	15.239	2045	2049	2055	rVB2	109191	159240	1.19%	0.096%
66	15.345	2063	2067	2069	rBV	45637	69642	0.52%	0.042%
67	15.386	2069	2074	2080	rVV	1238210	1675564	12.54%	1.005%
68	15.463	2080	2087	2094	rVB	2304815	3030735	22.68%	1.818%
69	15.645	2111	2118	2130	rVB	6552066	9377249	70.17%	5.626%
70	15.751	2130	2136	2146	rBV	1086325	1465772	10.97%	0.879%
71	16.004	2172	2179	2187	rBV2	950730	1399494	10.47%	0.840%
72	16.098	2192	2195	2200	rVB6	46493	67364	0.50%	0.040%
73	16.257	2218	2222	2224	rBV	74983	92114	0.69%	0.055%
74	16.292	2224	2228	2232	rVV2	75058	100770	0.75%	0.060%
75	16.345	2232	2237	2242	rVB	121456	169094	1.27%	0.101%
76	16.439	2250	2253	2257	rVB	61230	79542	0.60%	0.048%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012811.D
Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

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	16.651	2284	2289	2295	rBV	284840	413529	3.09%	0.248%
78	16.763	2301	2308	2314	rBV4	148197	332608	2.49%	0.200%
79	16.863	2320	2325	2328	rBV4	39032	66403	0.50%	0.040%
80	17.157	2372	2375	2379	rVV	172295	204915	1.53%	0.123%
81	17.198	2379	2382	2388	rVB2	90798	127596	0.95%	0.077%
82	17.263	2389	2393	2402	rVB3	64101	95766	0.72%	0.057%
83	17.357	2404	2409	2412	rBV5	48869	82304	0.62%	0.049%
84	17.410	2412	2418	2427	rVB	5162705	7394017	55.33%	4.436%
85	17.504	2428	2434	2446	rBV	7616006	10836200	81.09%	6.501%
86	17.686	2462	2465	2473	rVV	145489	214855	1.61%	0.129%
87	17.763	2475	2478	2484	rVB2	39023	65602	0.49%	0.039%
88	18.069	2525	2530	2534	rBV	155112	194749	1.46%	0.117%
89	18.269	2555	2564	2569	rBV6	129852	271097	2.03%	0.163%
90	18.357	2573	2579	2596	rVB	7516721	10504682	78.61%	6.302%
91	18.645	2624	2628	2632	rBV5	86012	112525	0.84%	0.068%
92	18.851	2659	2663	2669	rVB3	63524	88817	0.66%	0.053%
93	19.310	2737	2741	2747	rVB2	91155	119702	0.90%	0.072%
94	19.374	2748	2752	2757	rBV	98282	138112	1.03%	0.083%
95	19.439	2758	2763	2770	rVB5	93969	152162	1.14%	0.091%
96	19.504	2771	2774	2778	rBV2	81876	94140	0.70%	0.056%
97	19.733	2807	2813	2826	rBV	10224812	13363442	100.00%	8.017%
98	21.492	3107	3112	3124	rBV	6221015	8520889	63.76%	5.112%
99	23.851	3505	3513	3520	rBV2	6370131	12735822	95.30%	7.641%
100	24.010	3533	3540	3554	rVB2	4009730	8503934	63.64%	5.102%

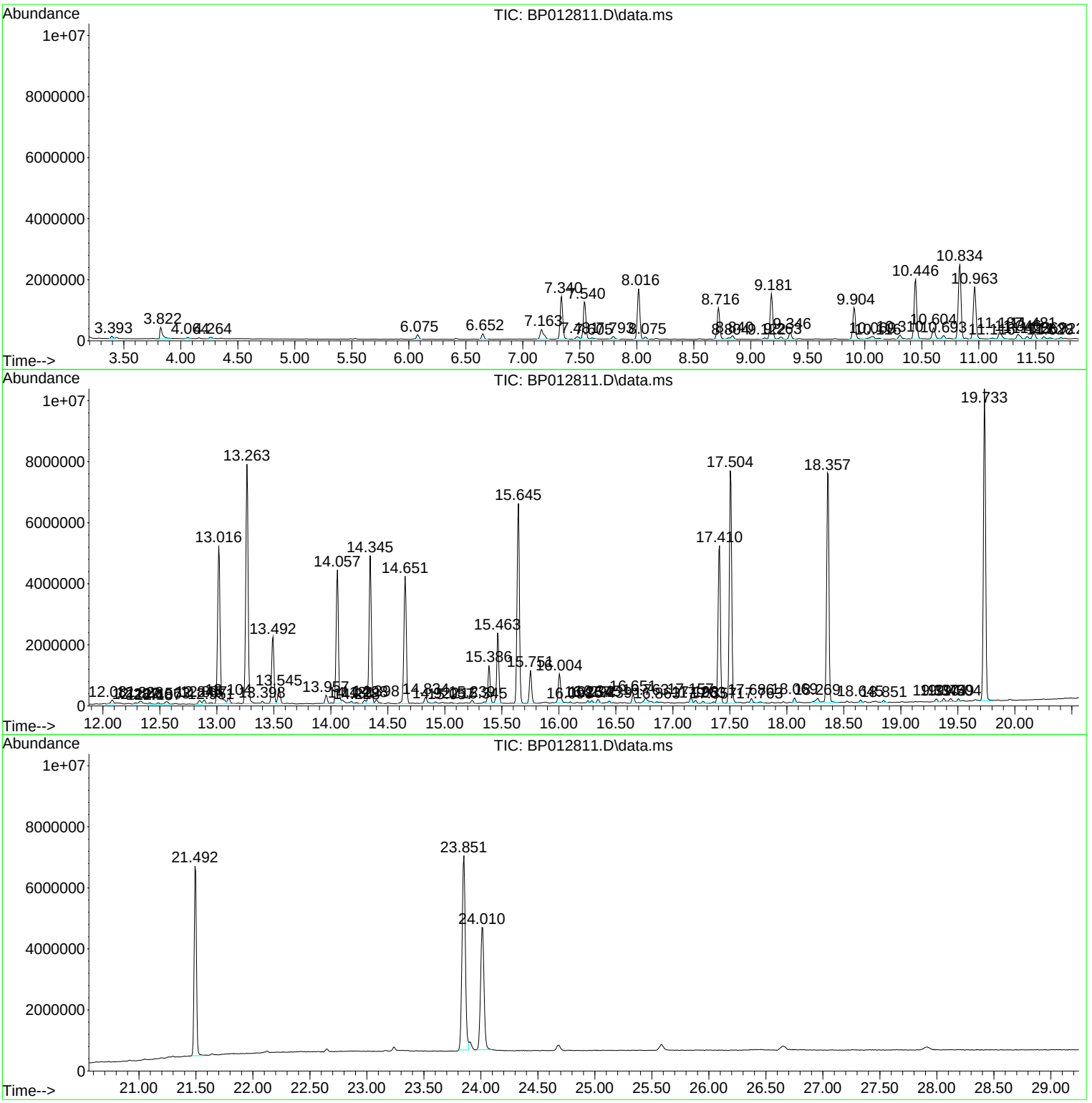
Sum of corrected areas: 166688039

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012811.D
Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

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 : BP012811.D
Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

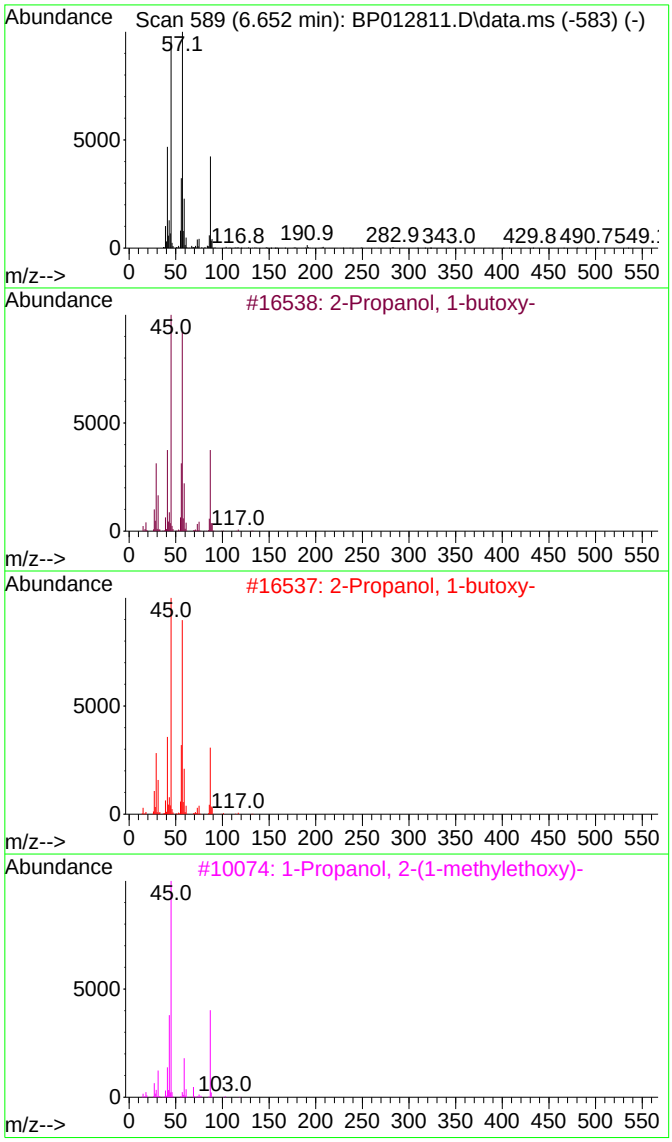
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

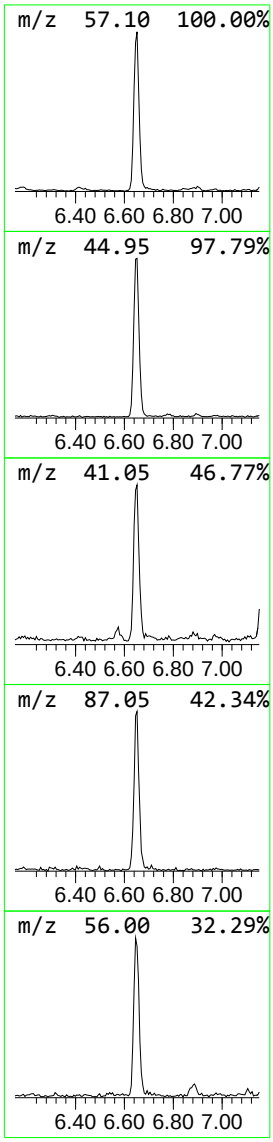
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.652	2.11 ng/ul	273618	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	90
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	Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012811.D
Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

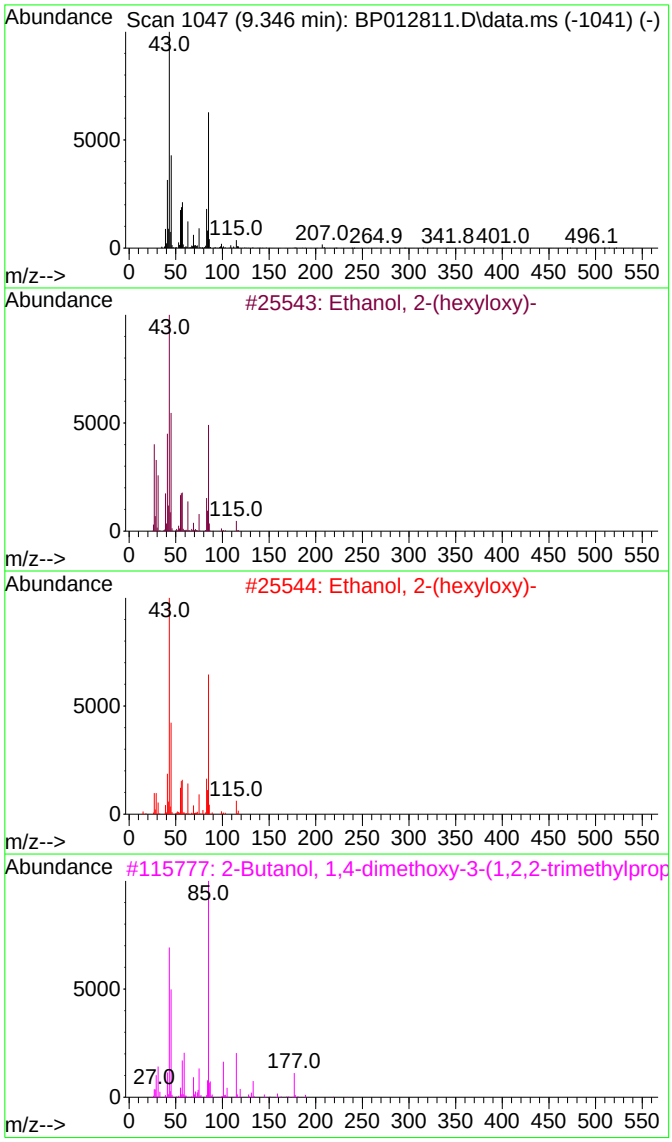
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

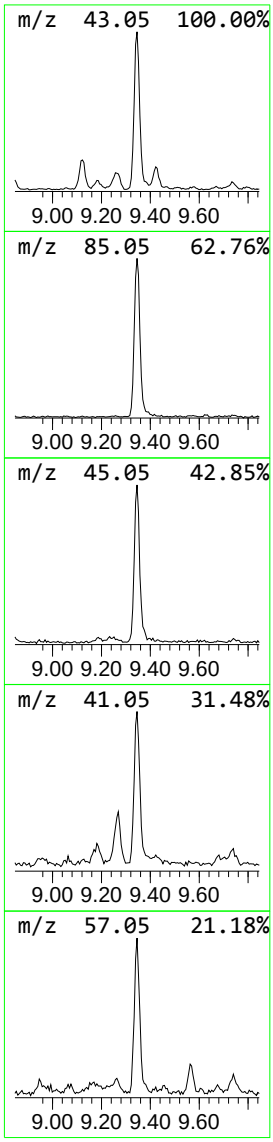
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.346	2.85 ng/ul	369274	1,4-Dichlorobenzene-d4	8.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	78
2	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	74
3	2-Butanol, 1,4-dimethoxy-3-(1,2,...	234	C12H26O4	079449-92-6	59
4	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	50
5	1-Bromo-8-tetrahydropyranyloxyoc...	292	C13H25BrO2	050816-20-1	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012811.D
Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

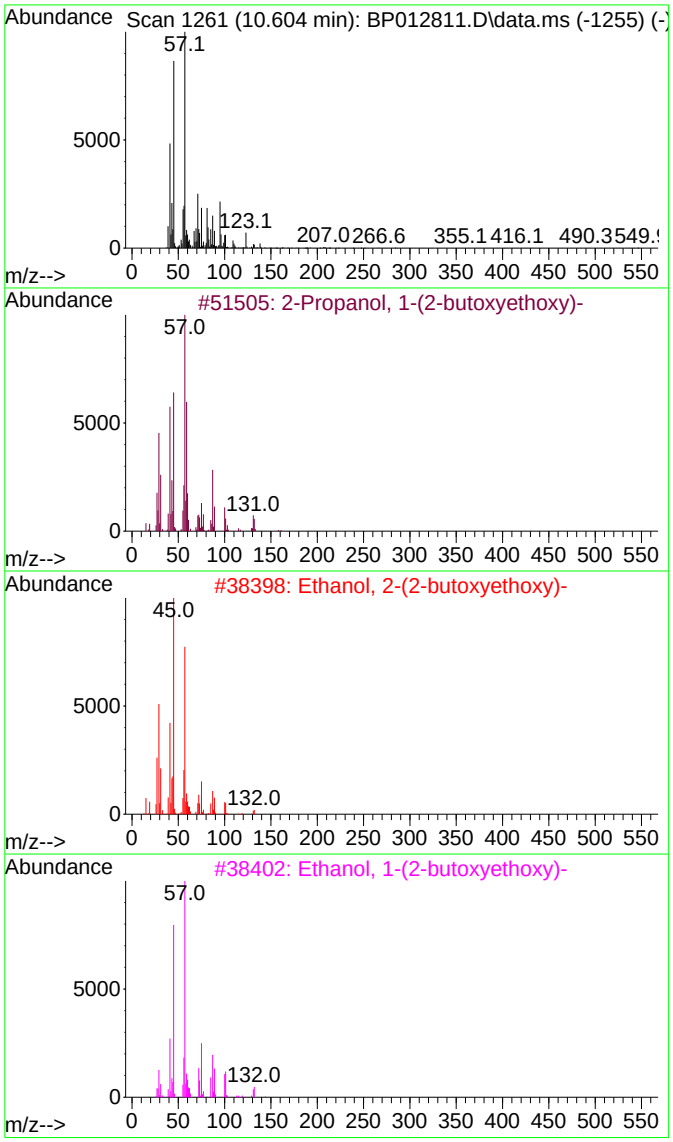
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

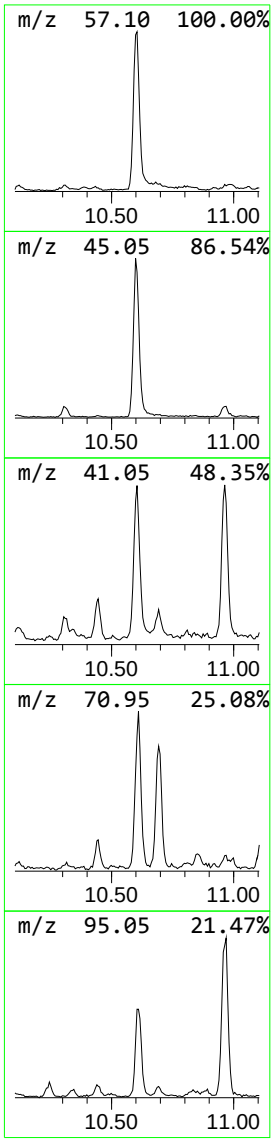
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Propanol, 1-(2-butoxyetho... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.604	3.21 ng/ul	662482	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-butoxyethoxy)-	176	C9H20O3	000124-16-3	53
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	50
3	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	49
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	49
5	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012811.D
Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

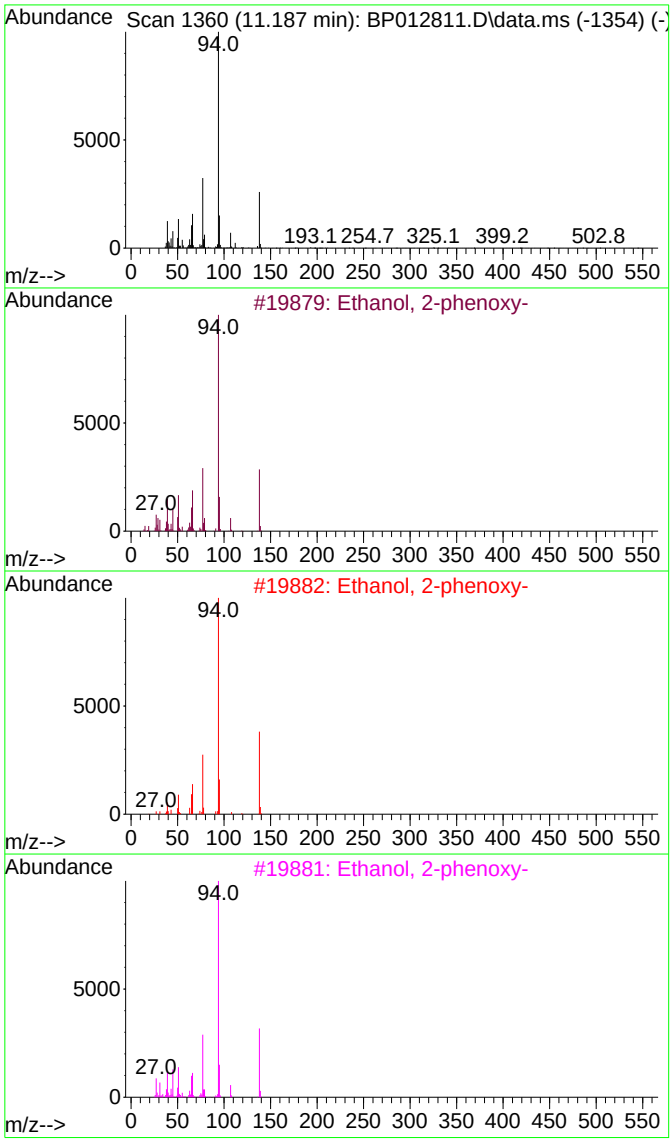
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

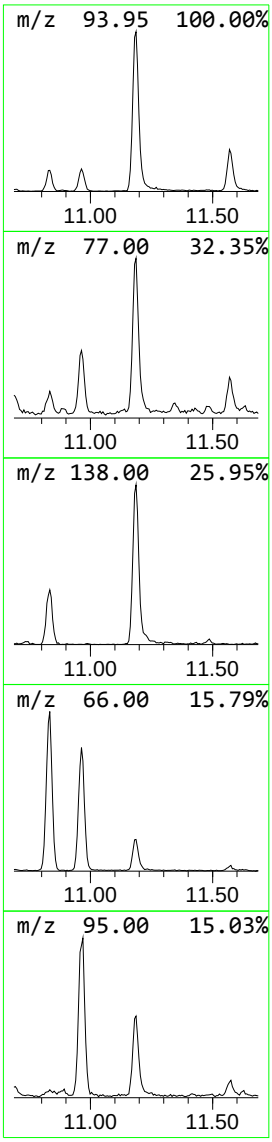
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Ethanol, 2-phenoxy- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.187	2.38 ng/ul	491176	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	97
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	91
5	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012811.D
Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

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

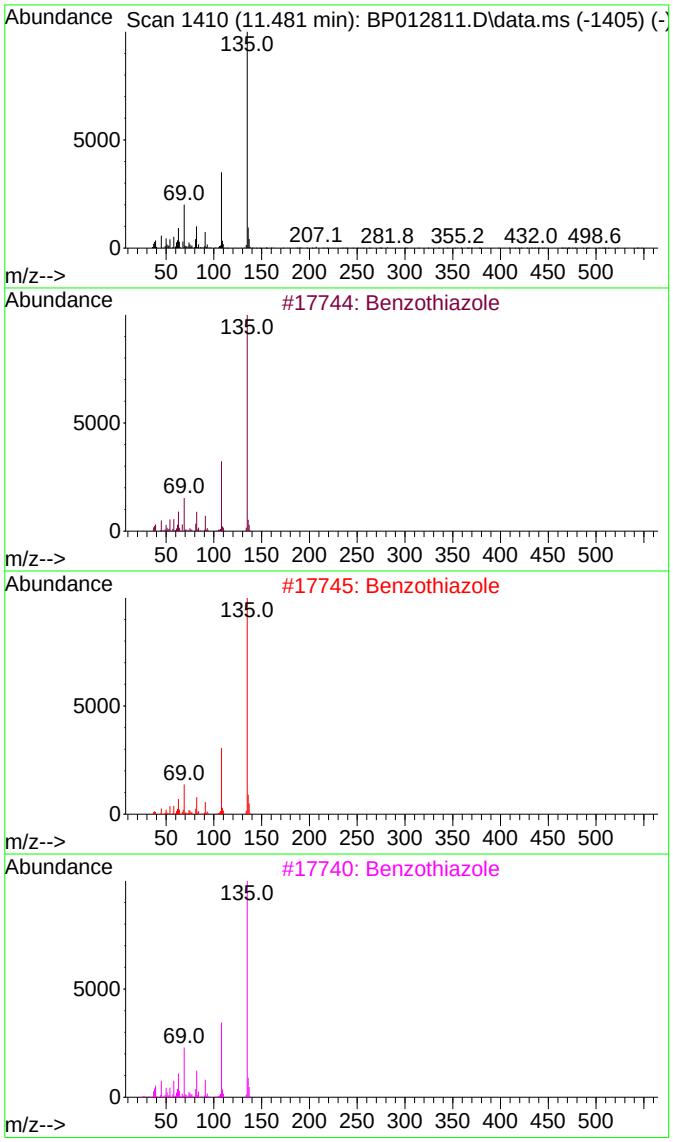
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

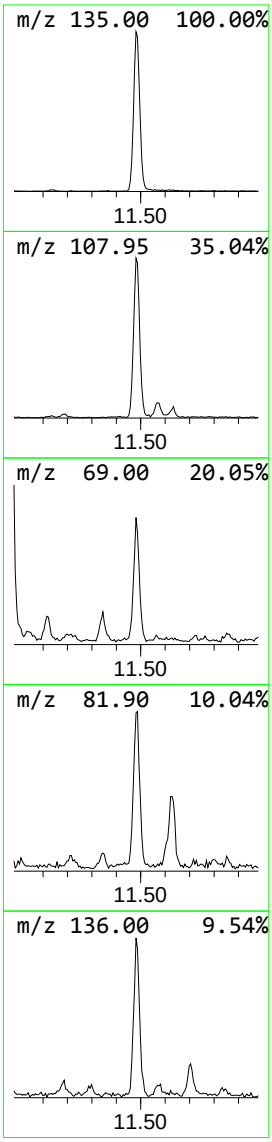
****

Peak Number 5 Benzothiazole Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.481	2.13 ng/ul	439249	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzothiazole	135	C7H5NS	000095-16-9	94
2	Benzothiazole	135	C7H5NS	000095-16-9	94
3	Benzothiazole	135	C7H5NS	000095-16-9	91
4	Benzothiazole	135	C7H5NS	000095-16-9	87
5	Benzothiazole	135	C7H5NS	000095-16-9	87





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Data File : BP012811.D
Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

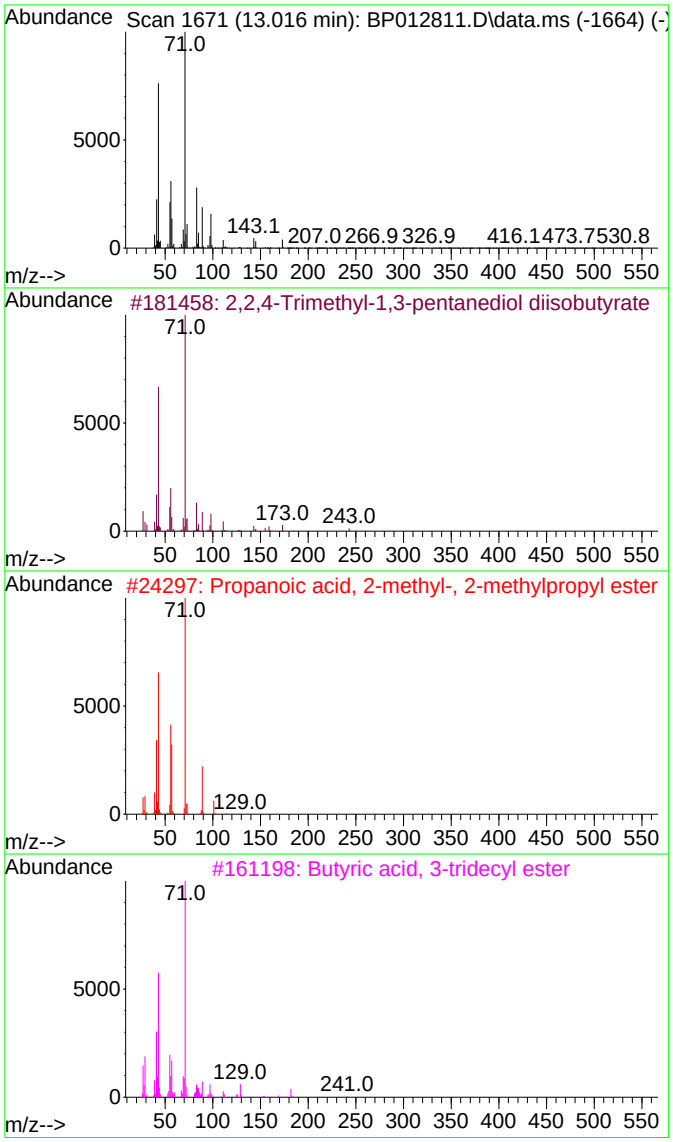
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

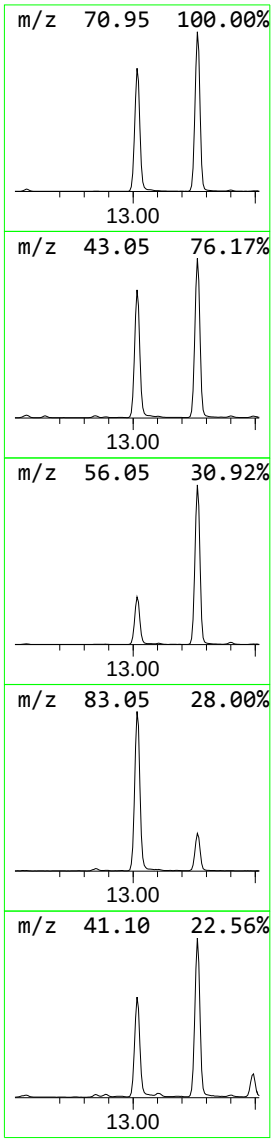
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 2,2,4-Trimethyl-1,3-pentane... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.016	25.58 ng/ul	7855370	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	Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	50
3	Butyric acid, 3-tridecyl ester	270	C17H34O2	1000280-56-8	47
4	Isophytol	296	C20H40O	000505-32-8	43
5	Butanoic acid, 1-methylpropyl ester	144	C8H16O2	000819-97-6	42





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Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

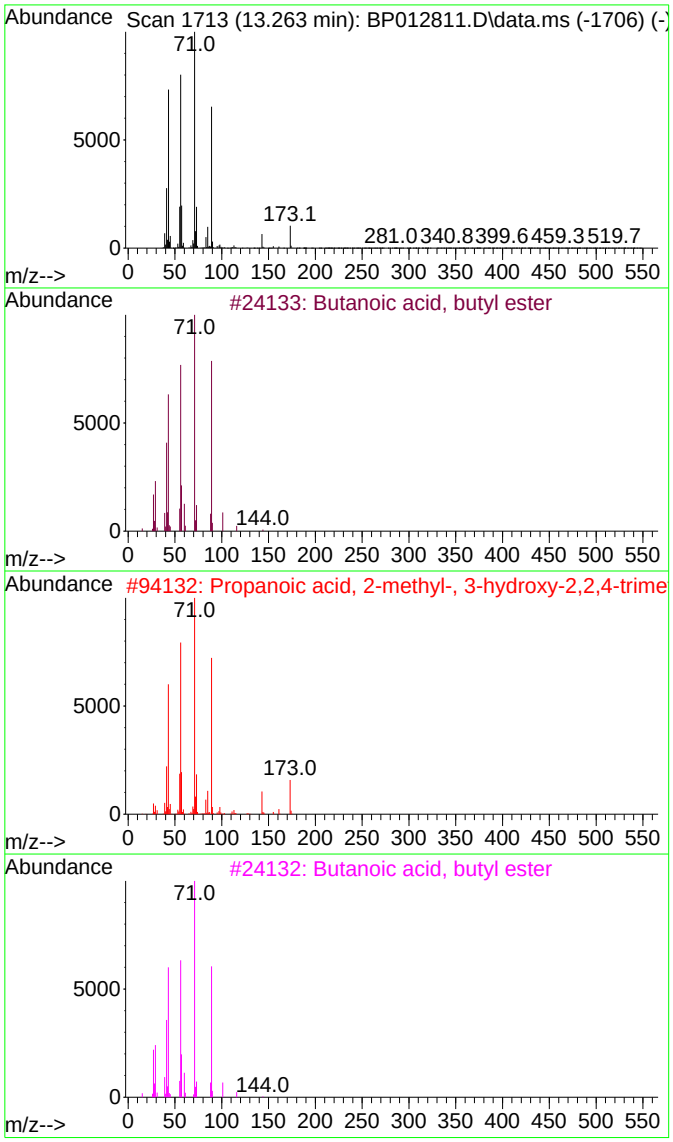
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

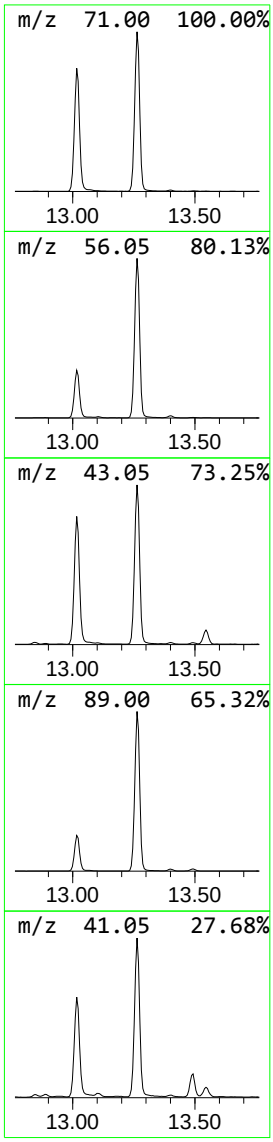
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Butanoic acid, butyl ester Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.263	36.45 ng/ul	11195900	Acenaphthene-d10	14.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	87
2	Propanoic acid, 2-methyl-, 3-hyd...	216	C12H24O3	000077-68-9	86
3	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	64
4	Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	64
5	Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	53





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Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

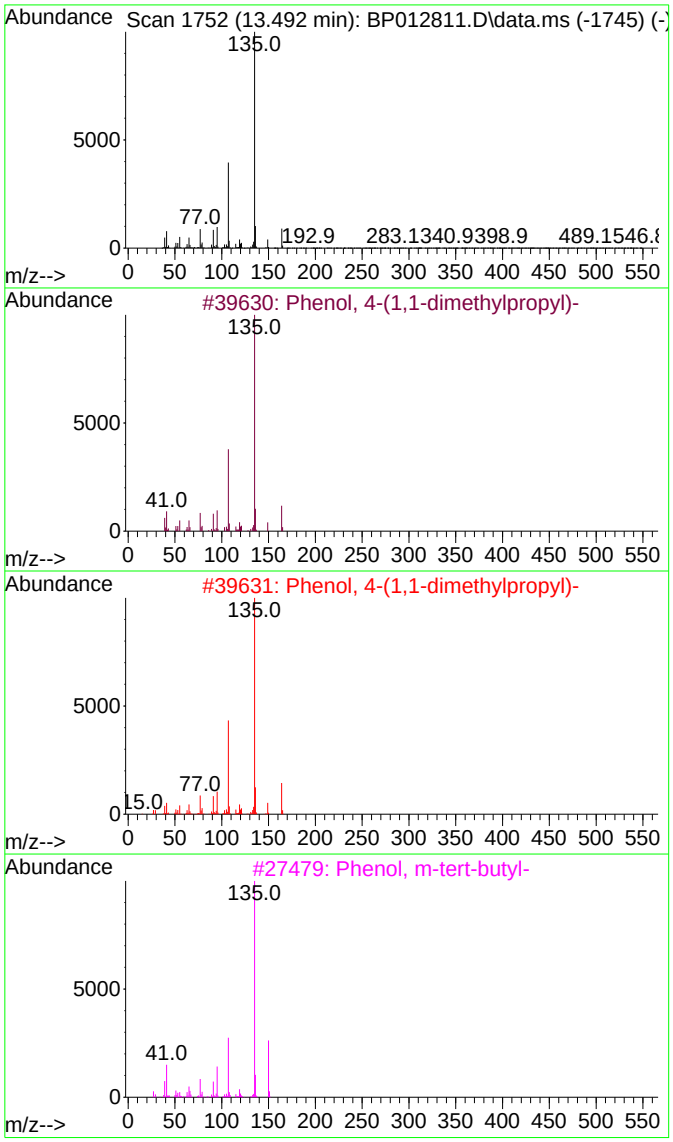
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

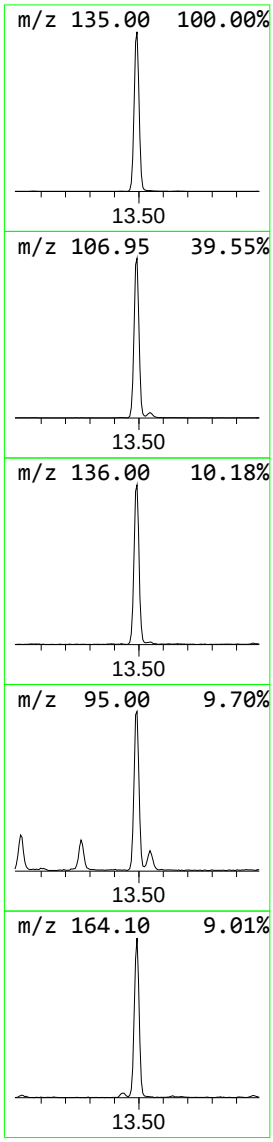
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Phenol, 4-(1,1-dimethylprop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.492	10.01 ng/ul	3074630	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, m-tert-butyl-	150	C10H14O	000585-34-2	86
4	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	81
5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	80





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Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

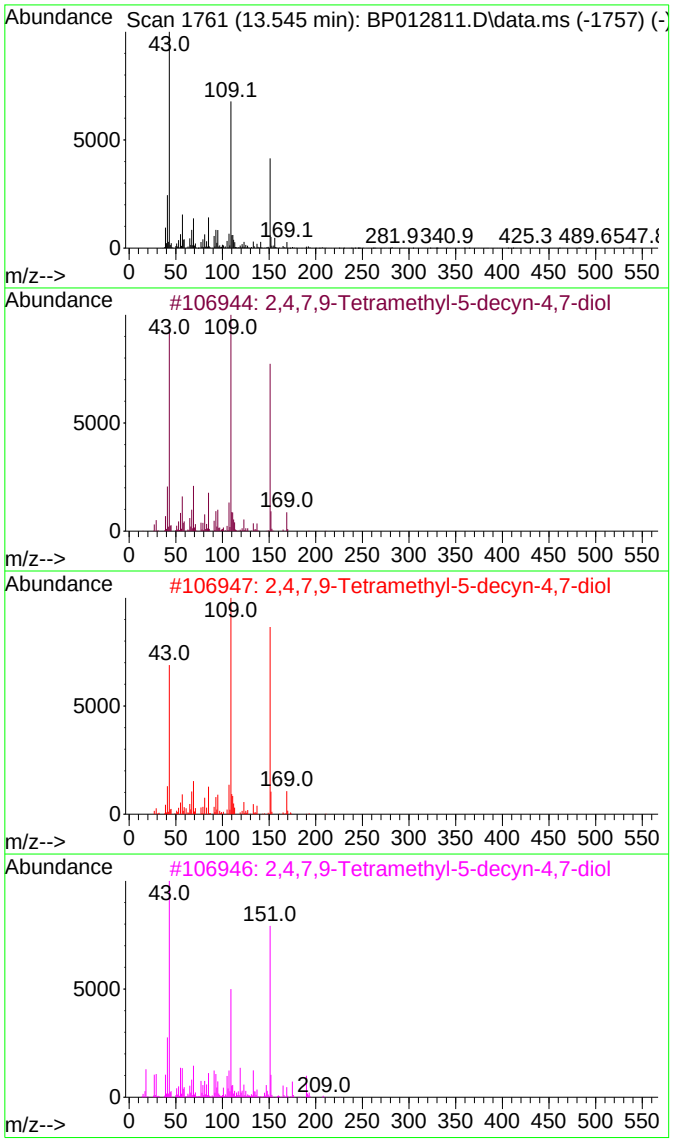
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

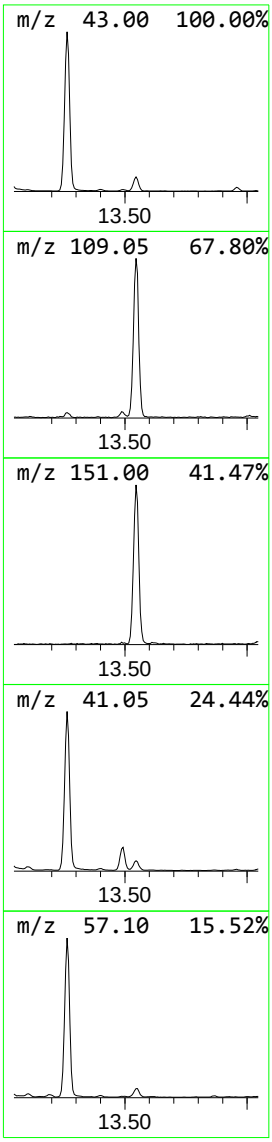
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.545	2.41 ng/ul	739139	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	86
3	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	58
4	m-Isopropoxyaniline	151	C9H13NO	041406-00-2	58
5	p-Isopropoxyaniline	151	C9H13NO	007664-66-6	58





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Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

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

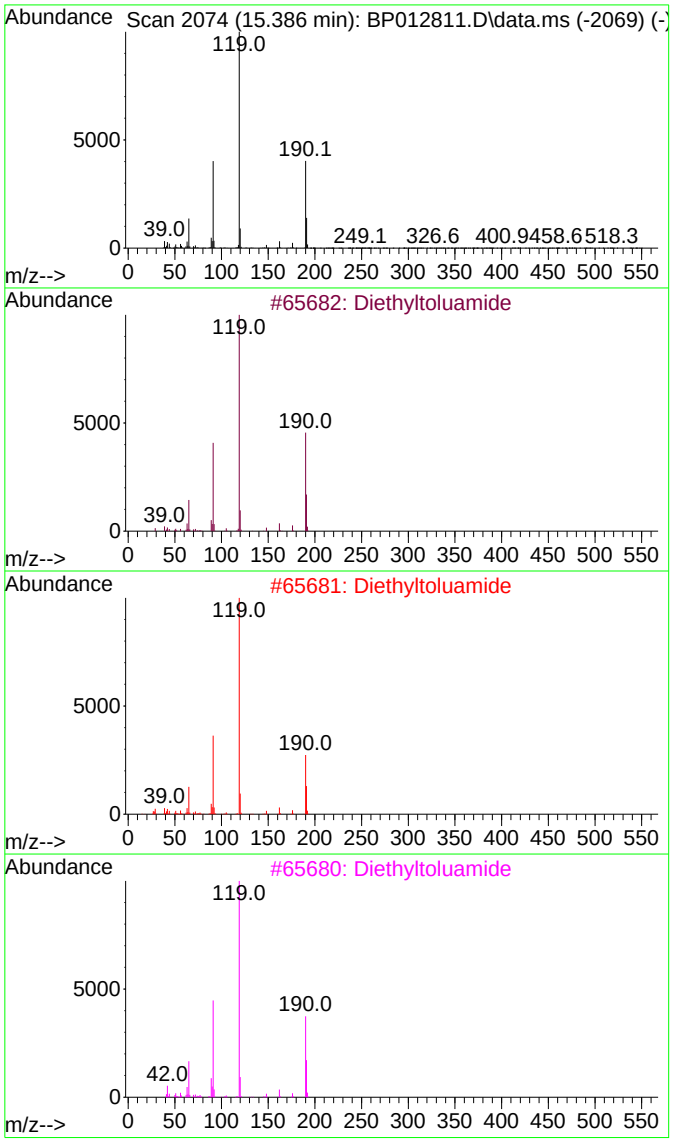
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

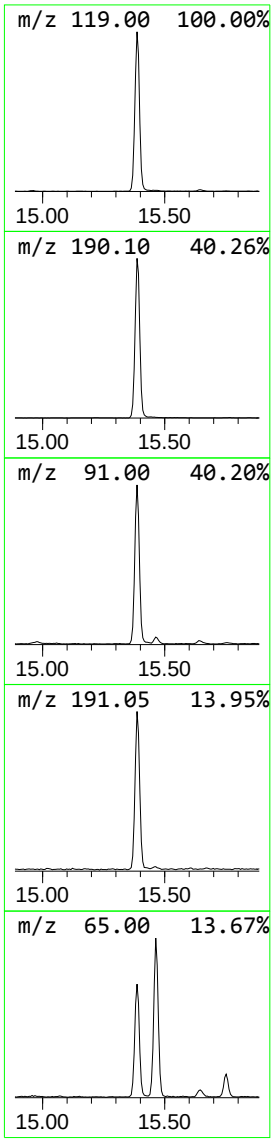
****

Peak Number 10 Diethyltoluamide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.387	5.46 ng/ul	1675560	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	Diethyltoluamide	191	C12H17NO	000134-62-3	91
4	Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	87
5	Benzamide, 3-methyl-N-butyl-	191	C12H17NO	1000407-41-1	76





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Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

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

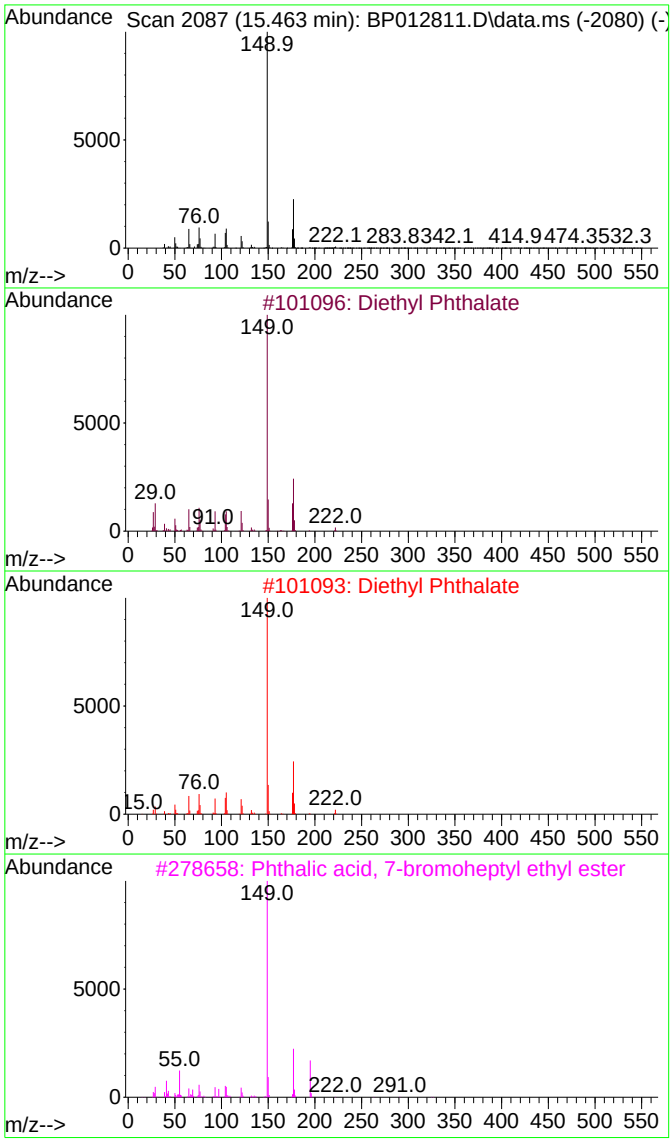
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

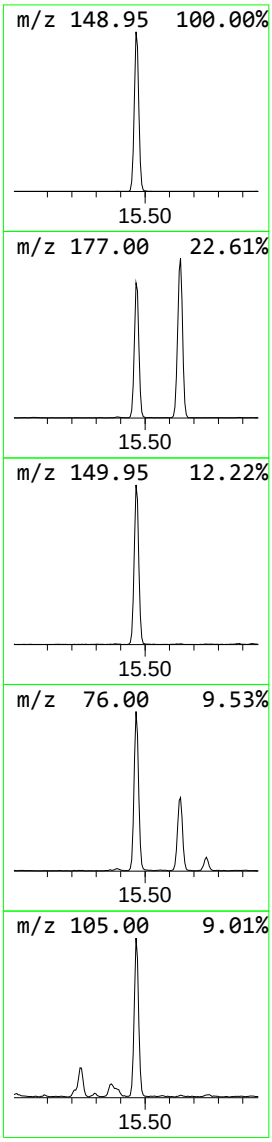
****

Peak Number 11 Diethyl Phthalate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.463	9.87 ng/ul	3030740	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, 7-bromoheptyl eth...	370	C17H23BrO4	1000415-51-6	83
4	1,2-Benzenedicarboxylic acid, mo...	222	C12H14O4	000131-70-4	74
5	Phthalic acid, ethyl 2-methylpen...	278	C16H22O4	1000356-90-5	72





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Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

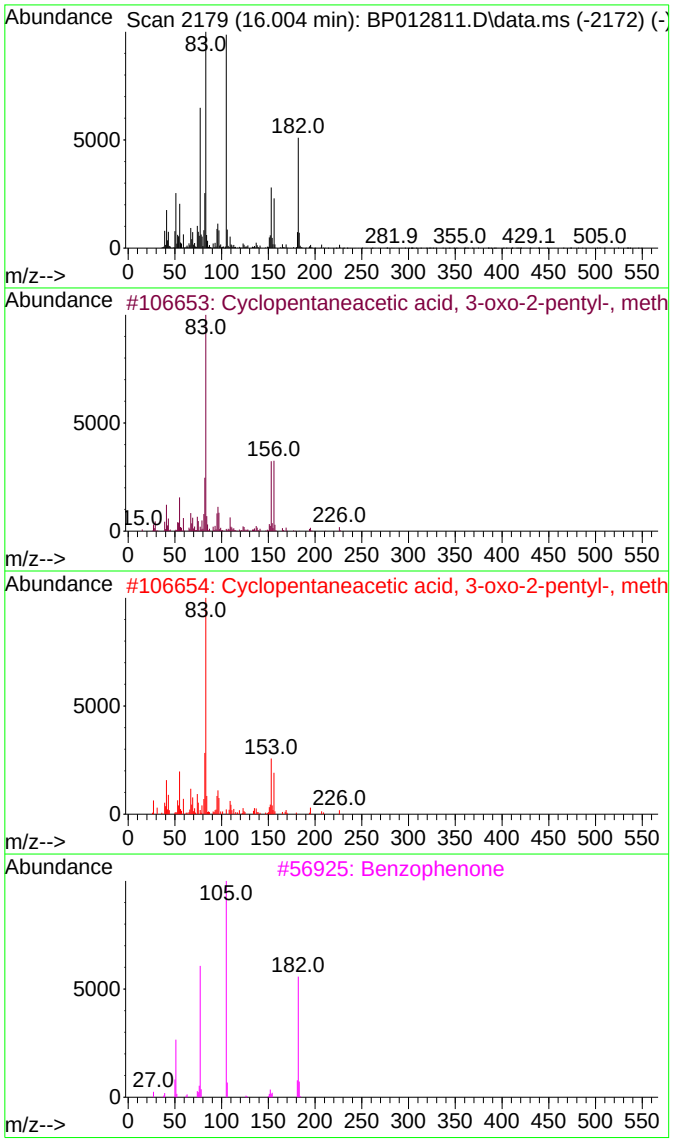
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

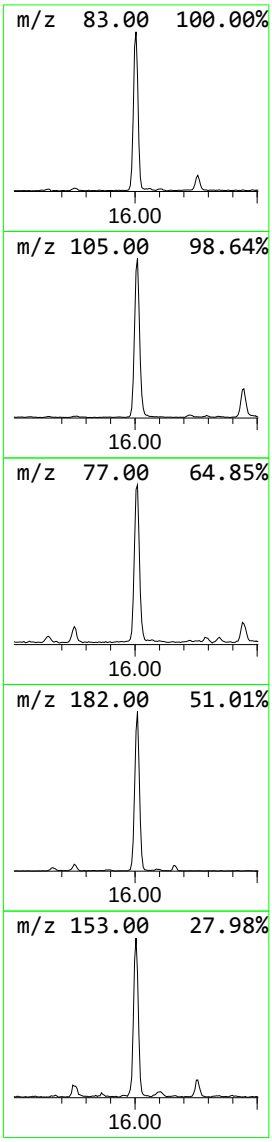
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Cyclopentaneacetic acid, 3-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.004	4.56 ng/ul	1399490	Acenaphthene-d10	14.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentaneacetic acid, 3-oxo-2...	226	C13H22O3	024851-98-7	78
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	42
5	Benzophenone	182	C13H10O	000119-61-9	41





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Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

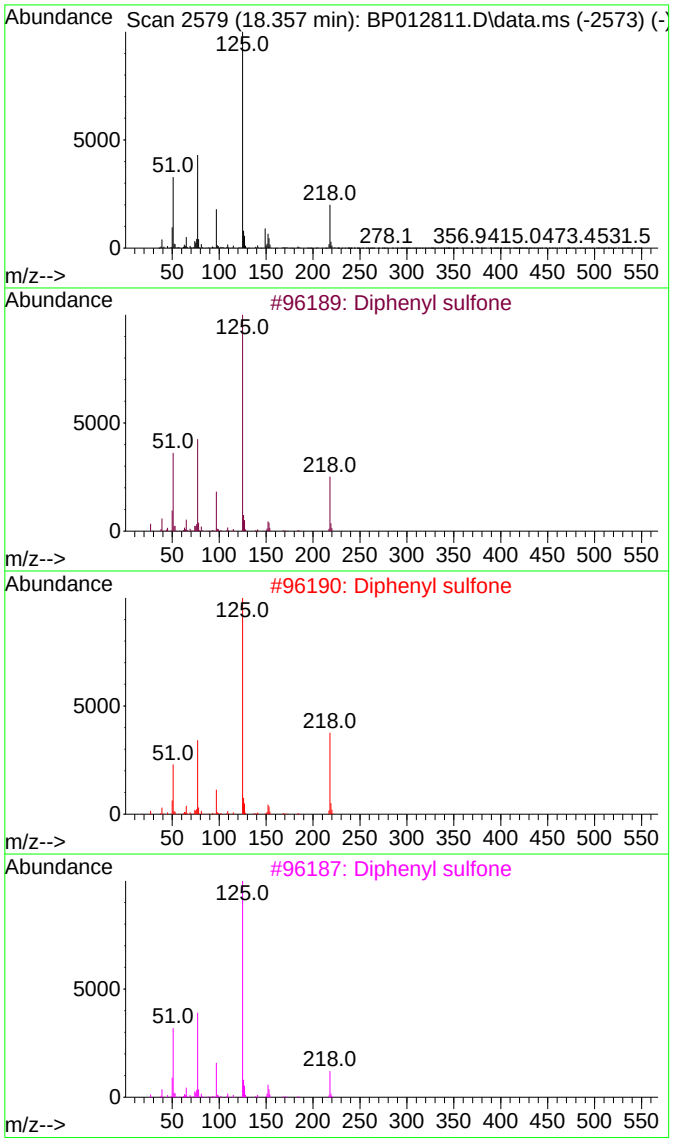
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

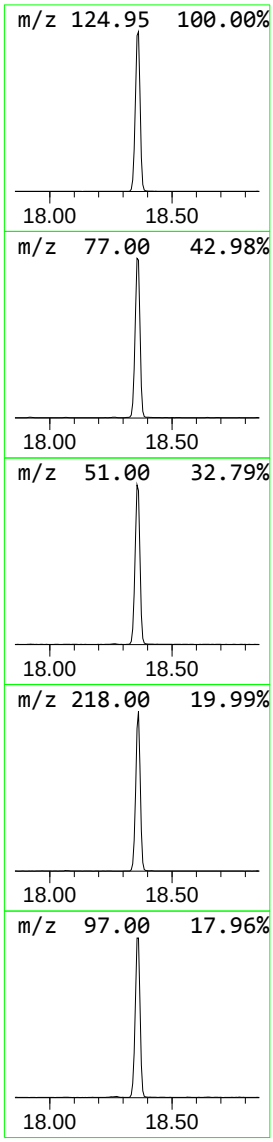
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Diphenyl sulfone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.357	28.41 ng/ul	10504700	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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012811.D
Acq On : 28 Nov 2022 11:53
Operator : CG/JU
Sample : N5710-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KY4

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
2-Propanol, 1-b...	6.652	2.1	ng/ul	273618	1	8.016	2591520	20.0
Ethanol, 2-(hex...	9.346	2.9	ng/ul	369274	1	8.016	2591520	20.0
2-Propanol, 1-(...	10.604	3.2	ng/ul	662482	2	10.834	4128540	20.0
Ethanol, 2-phen...	11.187	2.4	ng/ul	491176	2	10.834	4128540	20.0
Benzothiazole	11.481	2.1	ng/ul	439249	2	10.834	4128540	20.0
2,2,4-Trimethyl...	13.016	25.6	ng/ul	7855370	3	14.651	6142310	20.0
Butanoic acid, ...	13.263	36.5	ng/ul	11195900	3	14.651	6142310	20.0
Phenol, 4-(1,1-...	13.492	10.0	ng/ul	3074630	3	14.651	6142310	20.0
2,4,7,9-Tetrame...	13.545	2.4	ng/ul	739139	3	14.651	6142310	20.0
Diethyltoluamide	15.387	5.5	ng/ul	1675560	3	14.651	6142310	20.0
Diethyl Phthalate	15.463	9.9	ng/ul	3030740	3	14.651	6142310	20.0
Cyclopentaneace...	16.004	4.6	ng/ul	1399490	3	14.651	6142310	20.0
Diphenyl sulfone	18.357	28.4	ng/ul	10504700	4	17.410	7394020	20.0