LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040724\
Data File : BM045017.D
Acq On : 07 Apr 2024 20:55
Operator : MA/JU
Sample : P1970-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AL9

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_M\Methods\SFAM-EPA-BM040524.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BM045017.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.205	17	20	29	rVB	15559	23387	0.81%	0.083%
2	3.616	86	90	99	rBV	69102	128088	4.45%	0.454%
3	3.899	134	138	142	rBV5	6942	9237	0.32%	0.033%
4	4.151	177	181	183	rBV4	2745	4459	0.15%	0.016%
5	4.998	321	325	327	rBV5	2772	4145	0.14%	0.015%
6	5.728	445	449	450	rBV3	3781	3878	0.13%	0.014%
7	5.757	450	454	459	rVV3	5798	8643	0.30%	0.031%
8	5.928	477	483	486	rBB7	2790	4802	0.17%	0.017%
9	6.228	528	534	538	rBV8	3112	5014	0.17%	0.018%
10	6.663	603	608	612	rBV6	5696	7974	0.28%	0.028%
11	6.745	620	622	626	rVB5	2664	3551	0.12%	0.013%
12	6.816	626	634	646	rVB	64376	120188	4.18%	0.426%
13	6.981	656	662	672	rBV	317891	501684	17.44%	1.778%
14	7.163	686	693	704	rVV	303319	485455	16.87%	1.720%
15	7.628	765	772	778	rBV	307913	489191	17.00%	1.733%
16	7.692	778	783	791	rVV	64733	107048	3.72%	0.379%
17	7.816	802	804	812	rVB8	3515	6197	0.22%	0.022%
18	7.898	812	818	822	rBV6	8121	15172	0.53%	0.054%
19	8.222	868	873	875	rBV6	4680	7960	0.28%	0.028%
20	8.257	875	879	886	rVV4	8080	16703	0.58%	0.059%
21	8.339	886	893	901	rVV	211946	366319	12.73%	1.298%
22	8.404	903	904	908	rVV3	9204	9109	0.32%	0.032%
23	8.457	908	913	920	rVB2	16905	29891	1.04%	0.106%
24	8.722	953	958	963	rBV2	38167	61901	2.15%	0.219%
25	8.786	963	969	979	rVV	375649	651316	22.64%	2.308%
26	8.869	981	983	987	rVV4	6319	8506	0.30%	0.030%
27	8.904	987	989	992	rVV3	5186	6656	0.23%	0.024%
28	8.957	992	998	1006	rVB9	9627	21617	0.75%	0.077%
29	9.275	1045	1052	1054	rBV7	2585	4455	0.15%	0.016%
30	9.334	1058	1062	1065	rVB5	3694	4634	0.16%	0.016%
31	9.498	1084	1090	1104	rBV	301810	527047	18.32%	1.868%
32	9.692	1114	1123	1128	rBV10	10450	23210	0.81%	0.082%
33	9.751	1128	1133	1142	rVB7	11491	26790	0.93%	0.095%
34	9.928	1159	1163	1167	rVB8	2847	4455	0.15%	0.016%
35	10.033	1174	1181	1195	rBV	424490	768275	26.70%	2.722%
36	10.192	1204	1208	1218	rVB9	7712	19786	0.69%	0.070%

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Operator : MA/JU
Sample : P1970-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AL9

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_M\Methods\SFAM-EPA-BM040524.MA.M
Title : SVOA CALIBRATION

37	10.263	1218	1220	1224	rVB5	2362	3372	0.12%	0.012%
38	10.328	1224	1231	1237	rBV8	9339	23339	0.81%	0.083%
39	10.398	1237	1243	1250	rVV	420540	714593	24.84%	2.532%
40	10.469	1250	1255	1260	rVB6	12750	22965	0.80%	0.081%
41	10.557	1262	1270	1289	rBV	385214	727294	25.28%	2.577%
42	10.798	1305	1311	1313	rBV6	3456	4978	0.17%	0.018%
43	10.945	1329	1336	1341	rBV	559199	863659	30.02%	3.060%
44	11.004	1341	1346	1355	rVV	645653	988171	34.35%	3.502%
45	11.216	1373	1382	1389	rBV6	55473	117636	4.09%	0.417%
46	11.280	1389	1393	1399	rVV	17952	33636	1.17%	0.119%
47	11.345	1400	1404	1408	rVB5	3966	5911	0.21%	0.021%
48	11.775	1472	1477	1482	rVV8	2212	4597	0.16%	0.016%
49	11.933	1499	1504	1511	rVV9	3179	5309	0.18%	0.019%
50	12.004	1511	1516	1521	rVB4	7932	14733	0.51%	0.052%
51	12.169	1539	1544	1547	rBV5	4484	5794	0.20%	0.021%
52	12.363	1569	1577	1578	rBV6	2105	3704	0.13%	0.013%
53	12.486	1595	1598	1601	rBV4	2621	4256	0.15%	0.015%
54	12.563	1606	1611	1615	rVV7	9490	16943	0.59%	0.060%
55	12.616	1615	1620	1631	rVV	34160	63283	2.20%	0.224%
56	12.751	1638	1643	1651	rBV	6498	14510	0.50%	0.051%
57	12.874	1659	1664	1671	rBV	68206	104586	3.64%	0.371%
58	13.045	1689	1693	1698	rVB6	2729	5364	0.19%	0.019%
59	13.122	1698	1706	1708	rBV5	6585	9262	0.32%	0.033%
60	13.174	1708	1715	1726	rVV	634458	969499	33.70%	3.435%
61	13.245	1726	1727	1733	rVB5	3865	3998	0.14%	0.014%
62	13.704	1796	1805	1810	rBV	833561	1209289	42.03%	4.285%
63	13.751	1810	1813	1820	rVB2	14100	23390	0.81%	0.083%
64	13.869	1829	1833	1837	rBV6	2776	4401	0.15%	0.016%
65	13.969	1841	1850	1860	rBV	1069716	1576594	54.80%	5.587%
66	14.139	1875	1879	1882	rBV4	2189	3702	0.13%	0.013%
67	14.174	1882	1885	1891	rVV6	3361	7125	0.25%	0.025%
68	14.274	1895	1902	1912	rBV	687002	995897	34.61%	3.529%
69	14.545	1941	1948	1956	rBV5	19870	58637	2.04%	0.208%
70	14.639	1963	1964	1970	rVB6	3337	3567	0.12%	0.013%
71	14.716	1975	1977	1984	rVB4	3722	7045	0.24%	0.025%
72	14.886	2001	2006	2010	rBV6	4929	7374	0.26%	0.026%
73	14.974	2015	2021	2024	rBV6	5520	9470	0.33%	0.034%
74	15.039	2024	2032	2037	rVV	301191	407190	14.15%	1.443%
75	15.086	2037	2040	2042	rVV	8936	12404	0.43%	0.044%
76	15.121	2042	2046	2050	rVB	55976	71320	2.48%	0.253%

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Operator : MA/JU
Sample : P1970-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AL9

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_M\Methods\SFAM-EPA-BM040524.MA.M
Title : SVOA CALIBRATION

77	15.280	2066	2073	2083	rBV	1308427	1978245	68.76%	7.010%
78	15.421	2090	2097	2111	rBV	313892	494501	17.19%	1.752%
79	15.521	2111	2114	2118	rVB4	7846	8803	0.31%	0.031%
80	15.663	2132	2138	2144	rBV3	21285	29521	1.03%	0.105%
81	16.068	2203	2207	2210	rBV4	4048	5545	0.19%	0.020%
82	17.033	2365	2371	2380	rVB	813387	1161185	40.36%	4.115%
83	17.133	2381	2388	2404	rVB	1645900	2311895	80.36%	8.192%
84	17.333	2420	2422	2427	rBV3	4909	5480	0.19%	0.019%
85	17.715	2481	2487	2494	rBV	155447	201718	7.01%	0.715%
86	17.945	2523	2526	2531	rBV5	15035	21131	0.73%	0.075%
87	18.021	2536	2539	2545	rBV	10698	18607	0.65%	0.066%
88	18.151	2558	2561	2564	rBV3	8917	10516	0.37%	0.037%
89	18.239	2572	2576	2579	rVB4	3314	3787	0.13%	0.013%
90	18.415	2601	2606	2609	rBV8	3645	7382	0.26%	0.026%
91	18.527	2621	2625	2629	rVB7	4110	6647	0.23%	0.024%
92	18.756	2661	2664	2669	rVB6	2726	3685	0.13%	0.013%
93	19.003	2702	2706	2711	rBV3	20625	28347	0.99%	0.100%
94	19.080	2715	2719	2723	rVB2	14726	22525	0.78%	0.080%
95	19.239	2742	2746	2750	rBV5	13897	21150	0.74%	0.075%
96	19.439	2774	2780	2788	rBV	2065040	2738154	95.17%	9.703%
97	19.803	2840	2842	2844	rBV3	4404	4207	0.15%	0.015%
98	21.245	3082	3087	3094	rBV	946129	1260037	43.80%	4.465%
99	23.321	3433	3440	3447	rBV2	1622944	2877072	100.00%	10.195%
100	23.462	3458	3464	3475	rVB	729849	1384421	48.12%	4.906%

Sum of corrected areas: 28220101

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040724\
Data File : BM045017.D
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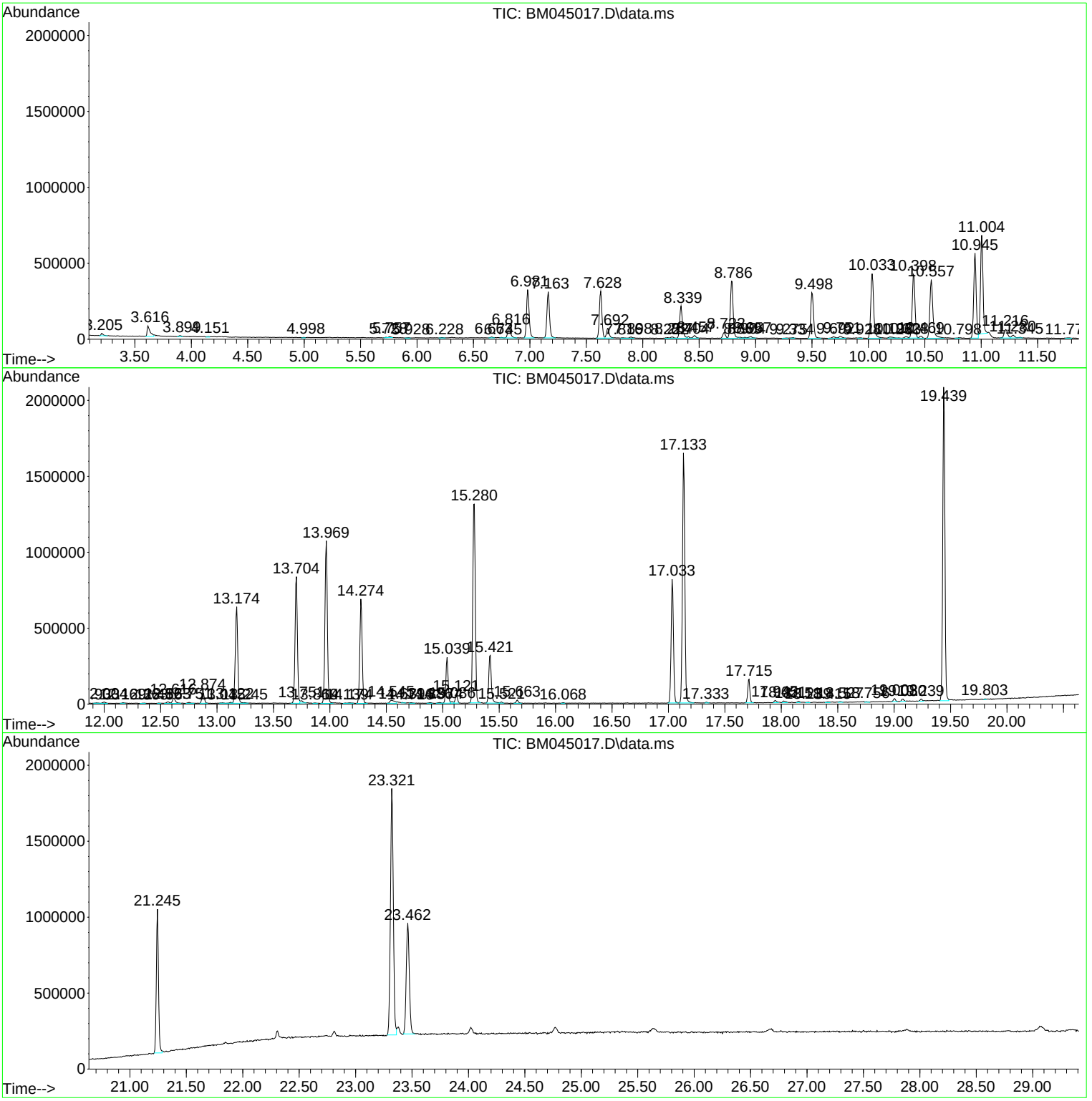
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AL9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040724\
Data File : BM045017.D
Acq On : 07 Apr 2024 20:55
Operator : MA/JU
Sample : P1970-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AL9

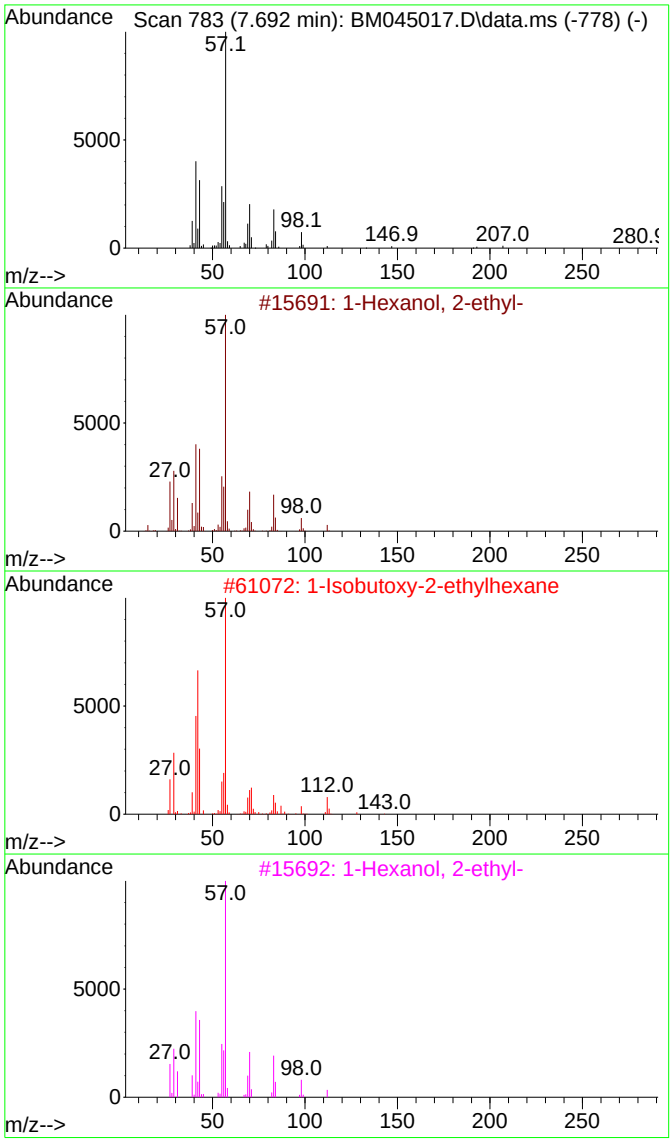
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

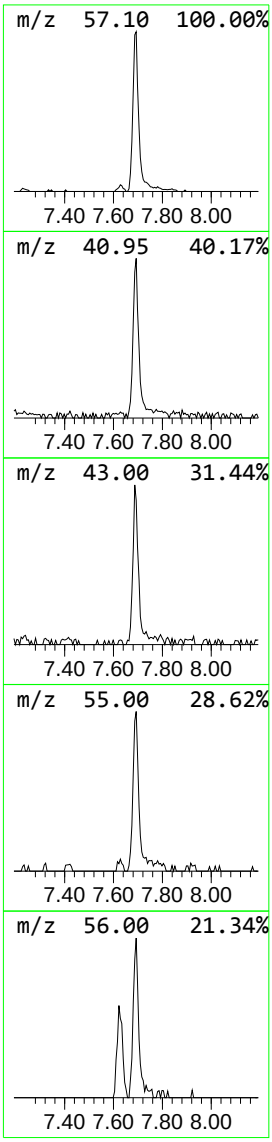
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.692	4.38 ng/ul	107048	1,4-Dichlorobenzene-d4	7.628

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2	1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	78
3	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	56





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040724\
Data File : BM045017.D
Acq On : 07 Apr 2024 20:55
Operator : MA/JU
Sample : P1970-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AL9

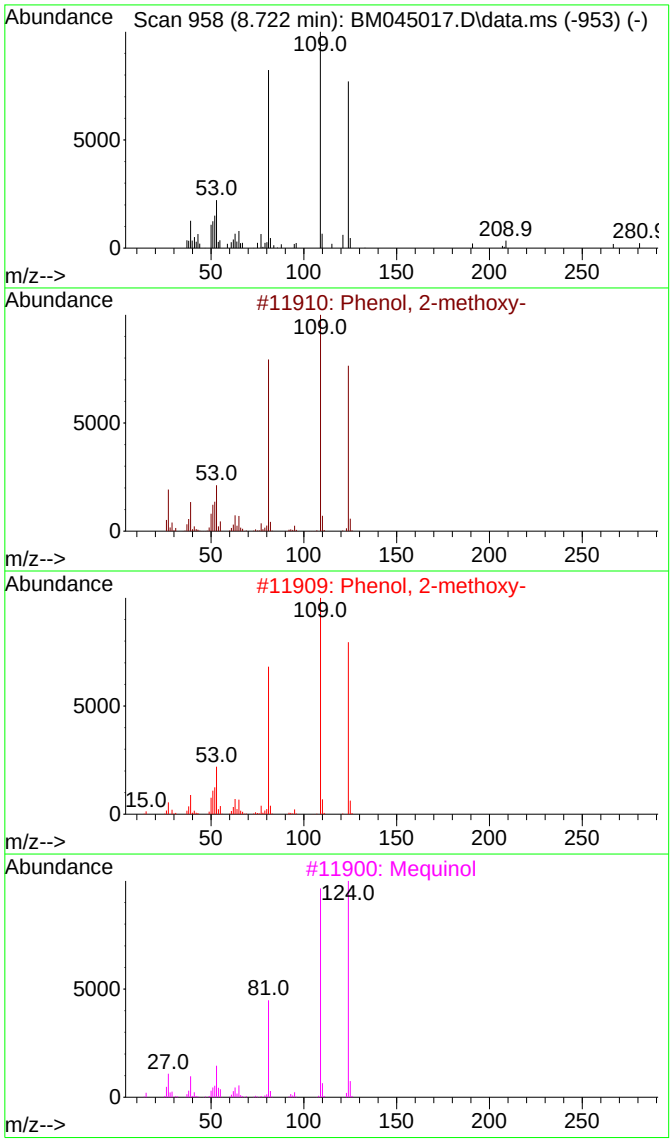
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

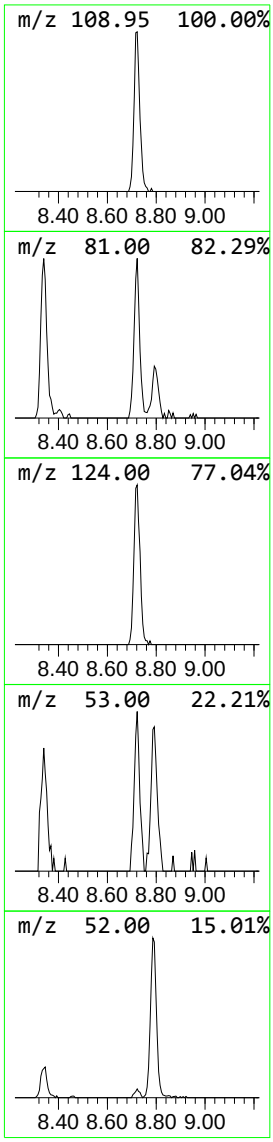
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Phenol, 2-methoxy- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.722	2.53 ng/ul	61901	1,4-Dichlorobenzene-d4	7.628

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	93
2	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	93
3	Mequinol	124	C7H8O2	000150-76-5	91
4	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	90
5	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	86





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ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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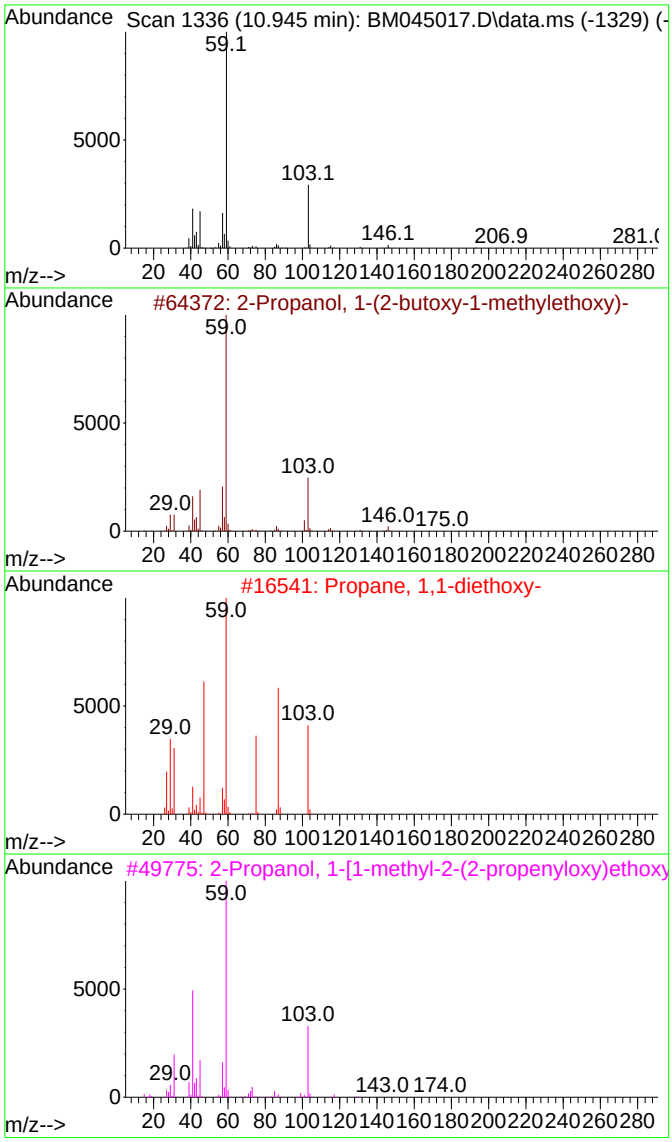
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

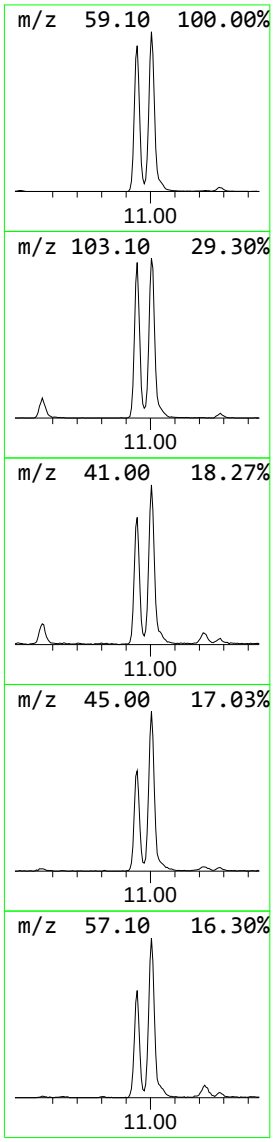
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Propanol, 1-(2-butoxy-1-m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.945	24.17 ng/ul	863659	Naphthalene-d8	10.398

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	83
2	Propane, 1,1-diethoxy-	132	C7H16O2	004744-08-5	78
3	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	74
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72
5	Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	64





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Instrument :
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ClientSampleId :
C0AL9

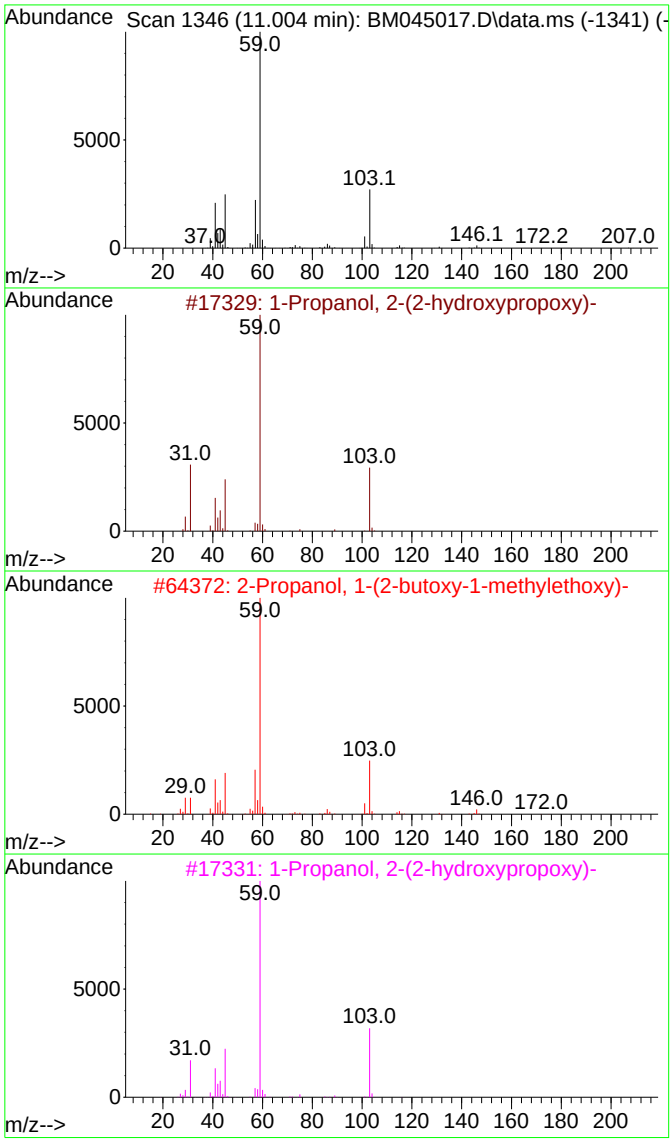
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

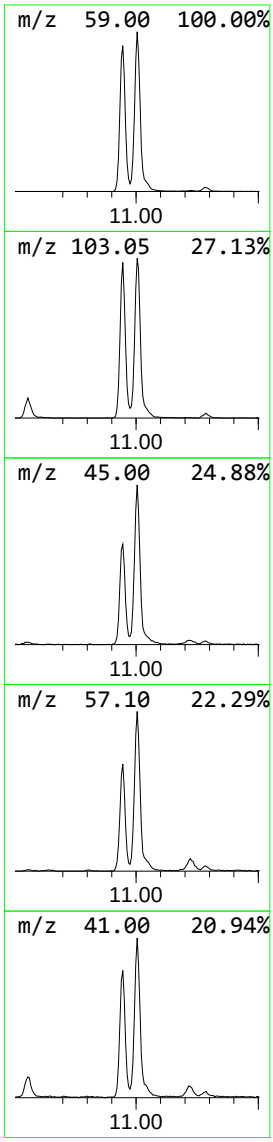
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1-Propanol, 2-(2-hydroxypro... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.004	27.66 ng/ul	988171	Naphthalene-d8	10.398

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78
2	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	78
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78
4	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	72
5	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	64





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Data File : BM045017.D
Acq On : 07 Apr 2024 20:55
Operator : MA/JU
Sample : P1970-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AL9

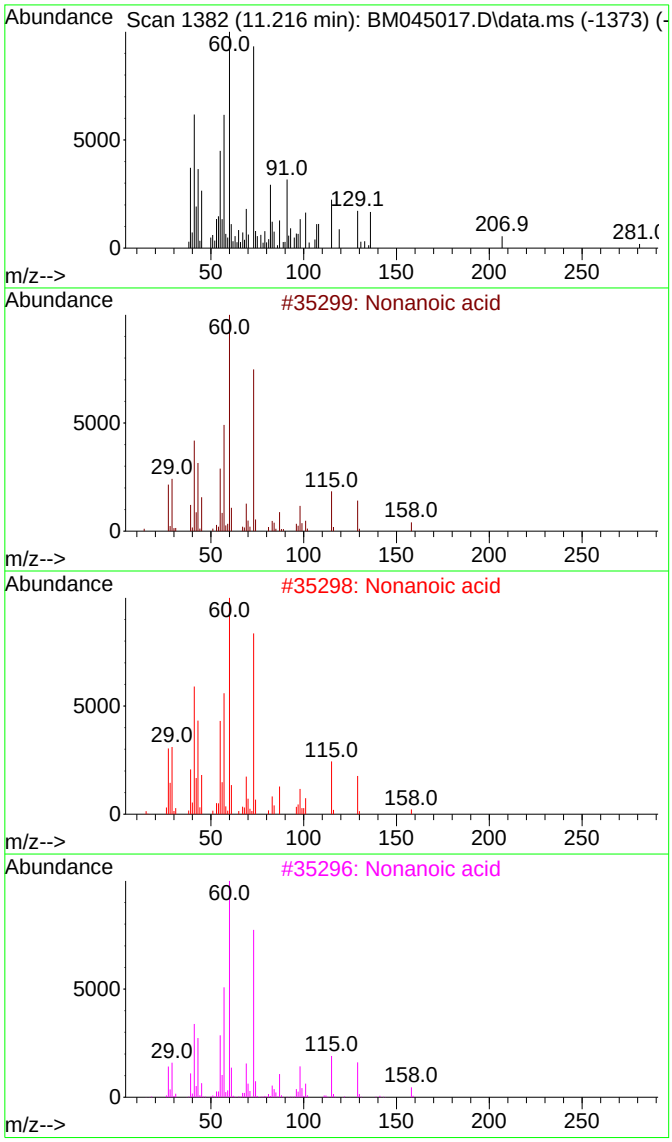
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

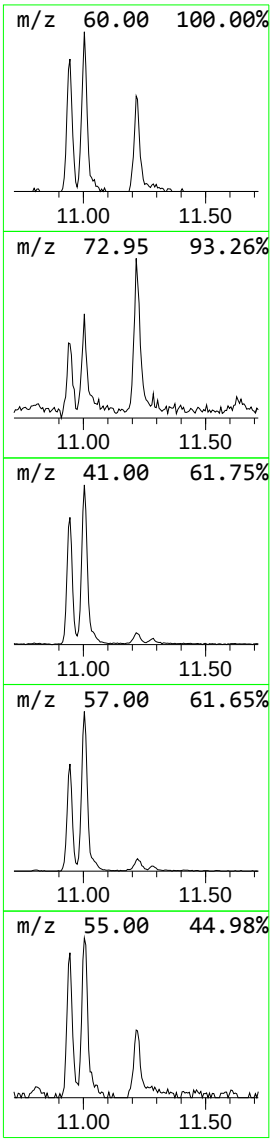
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Nonanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.216	3.29 ng/ul	117636	Naphthalene-d8	10.398

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid	158	C9H18O2	000112-05-0	72
2	Nonanoic acid	158	C9H18O2	000112-05-0	72
3	Nonanoic acid	158	C9H18O2	000112-05-0	72
4	Nonanoic acid	158	C9H18O2	000112-05-0	64
5	n-Decanoic acid	172	C10H20O2	000334-48-5	53





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Acq On : 07 Apr 2024 20:55
Operator : MA/JU
Sample : P1970-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AL9

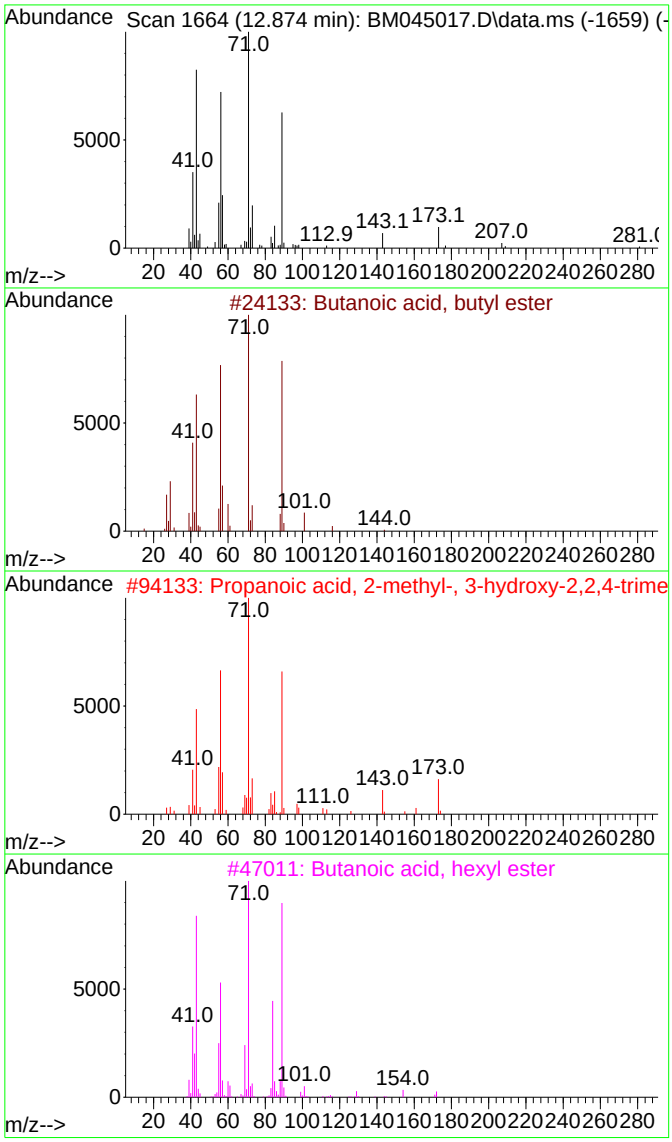
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

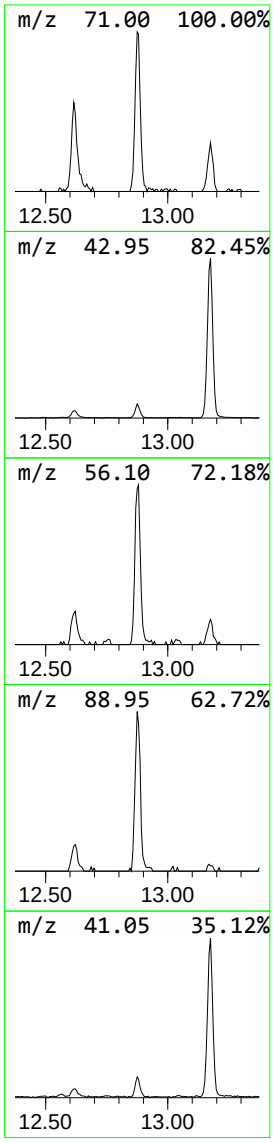
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Butanoic acid, butyl ester Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.874	2.10 ng/ul	104586	Acenaphthene-d10	14.274

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	64
2	Propanoic acid, 2-methyl-, 3-hyd...	216	C12H24O3	000077-68-9	64
3	Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	56
4	Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	56
5	Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	56





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Data File : BM045017.D
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Operator : MA/JU
Sample : P1970-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AL9

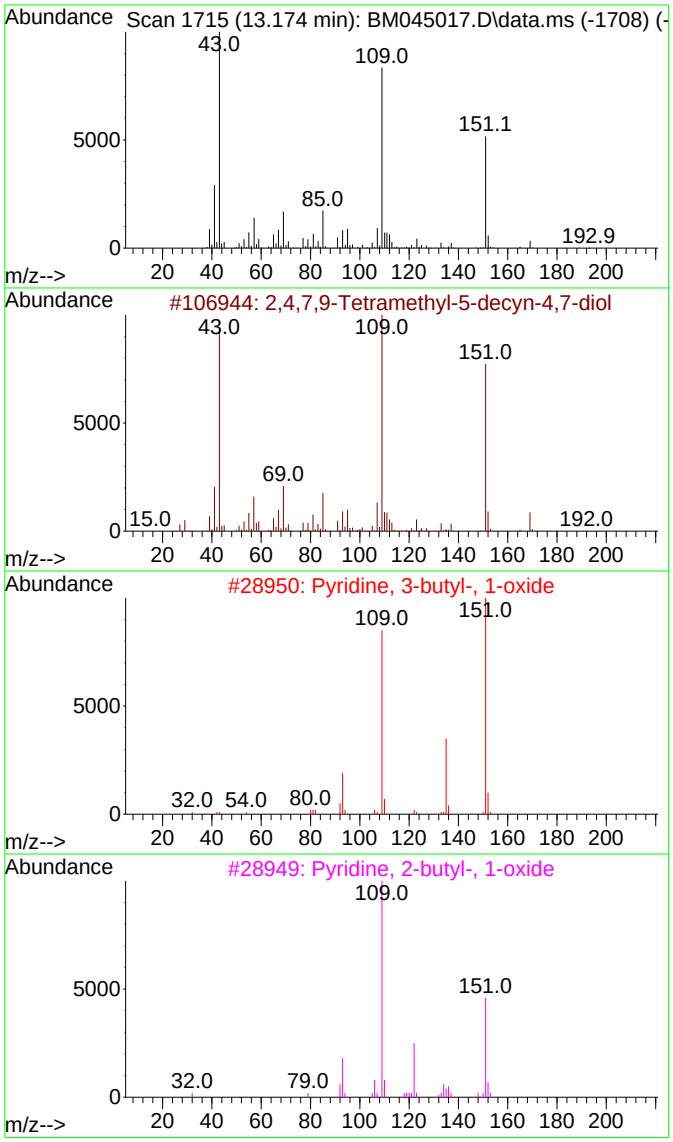
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

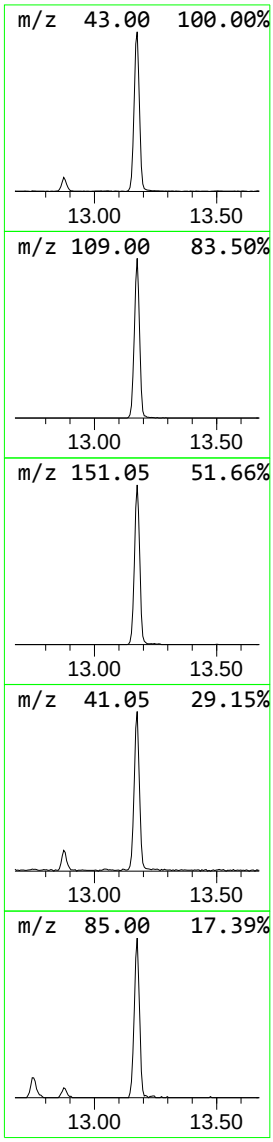
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.174	19.47 ng/ul	969499	Acenaphthene-d10	14.274

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	Pyridine, 3-butyl-, 1-oxide	151	C9H13NO	031396-33-5	59
3	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	59
4	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	59
5	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	58





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Data File : BM045017.D
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Operator : MA/JU
Sample : P1970-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AL9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

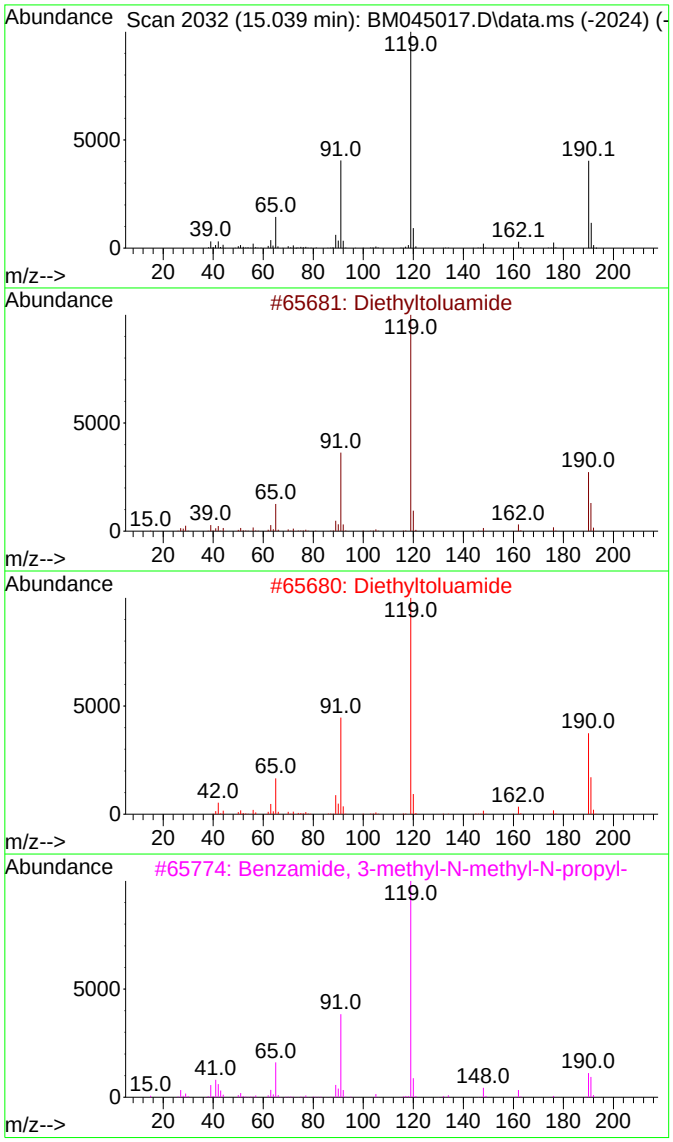
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

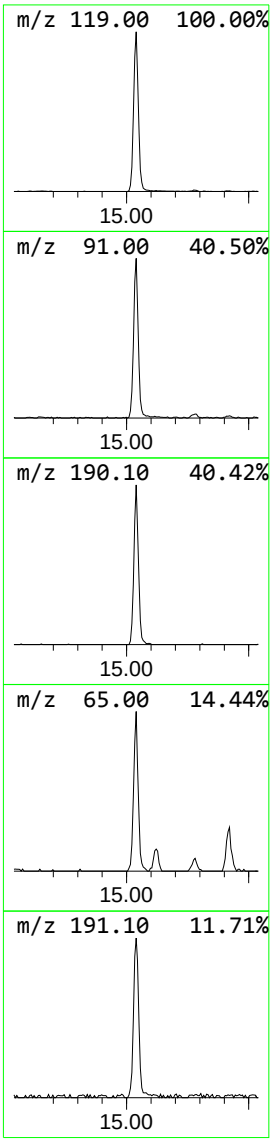
****

Peak Number 8 Diethyltoluamide Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.039	8.18 ng/ul	407190	Acenaphthene-d10	14.274

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	95
2	Diethyltoluamide	191	C12H17NO	000134-62-3	92
3	Benzamide, 3-methyl-N-methyl-N-p...	191	C12H17NO	1000421-46-2	81
4	L-Valine, N-(4-methylbenzoyl)-, ...	277	C16H23NO3	1000346-63-8	78
5	Benzamide, 3-methyl-N-butyl-	191	C12H17NO	1000407-41-1	76





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Data File : BM045017.D
Acq On : 07 Apr 2024 20:55
Operator : MA/JU
Sample : P1970-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AL9

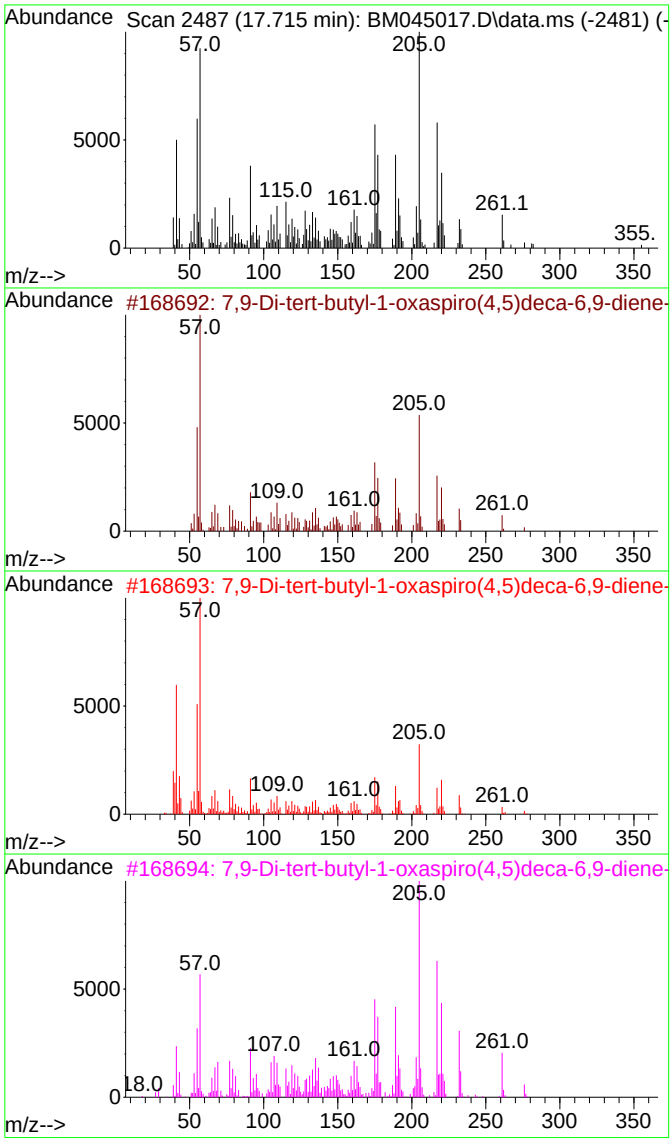
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION

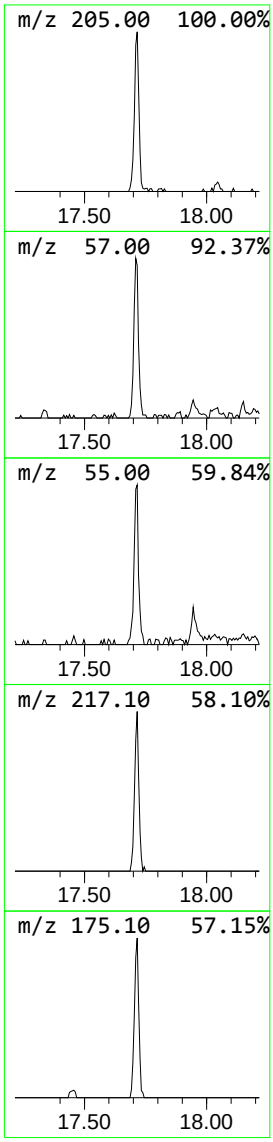
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 7,9-Di-tert-butyl-1-oxaspir... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.715	3.47 ng/ul	201718	Phenanthrene-d10	17.033

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	95
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	74
3	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	72
4	1-(3-Amino-4,6-dimethylthieno[2,...	220	C11H12N2OS	052505-42-7	50
5	Ethanone, 1-(5,6,7,8-tetrahydro-...	220	C14H20O2	071596-88-8	46





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040724\
Data File : BM045017.D
Acq On : 07 Apr 2024 20:55
Operator : MA/JU
Sample : P1970-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AL9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.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
1-Hexanol, 2-et...	7.692	4.4	ng/ul	107048	1	7.628	489191	20.0
Phenol, 2-methoxy-	8.722	2.5	ng/ul	61901	1	7.628	489191	20.0
2-Propanol, 1-(...	10.945	24.2	ng/ul	863659	2	10.398	714593	20.0
1-Propanol, 2-(...	11.004	27.7	ng/ul	988171	2	10.398	714593	20.0
Nonanoic acid	11.216	3.3	ng/ul	117636	2	10.398	714593	20.0
Butanoic acid, ...	12.874	2.1	ng/ul	104586	3	14.274	995897	20.0
2,4,7,9-Tetrame...	13.174	19.5	ng/ul	969499	3	14.274	995897	20.0
Diethyltoluamide	15.039	8.2	ng/ul	407190	3	14.274	995897	20.0
7,9-Di-tert-but...	17.715	3.5	ng/ul	201718	4	17.033	1161190	20.0