LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015117.D
Acq On : 26 Jun 2021 15:12
Operator : CG/JU
Sample : M2704-15
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDM5

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_N\METHODS\SFAM-EPA-BN061821MA.M
Title : SVOA CALIBRATION

Signal : TIC: BN015117.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.281	30	33	41	rVB	34341	47063	0.98%	0.113%
2	3.546	76	78	85	rVB2	8474	13618	0.28%	0.033%
3	3.675	97	100	111	rBV	272835	413578	8.64%	0.991%
4	3.840	125	128	136	rVB5	5563	11696	0.24%	0.028%
5	3.911	136	140	142	rVV4	5482	8751	0.18%	0.021%
6	3.999	150	155	161	rVB	57880	81433	1.70%	0.195%
7	4.358	211	216	222	rVB	20866	30787	0.64%	0.074%
8	4.487	233	238	241	rBV	17472	23410	0.49%	0.056%
9	4.534	241	246	254	rVB	58330	79004	1.65%	0.189%
10	5.281	365	373	380	rVB	40966	60397	1.26%	0.145%
11	5.411	389	395	404	rVB	126388	243814	5.09%	0.584%
12	5.769	451	456	461	rBV	58827	90007	1.88%	0.216%
13	5.840	464	468	476	rVB2	7480	12907	0.27%	0.031%
14	6.740	618	621	627	rVB4	6685	11537	0.24%	0.028%
15	6.893	639	647	658	rBV	544294	849777	17.76%	2.036%
16	7.063	667	676	688	rBV	417198	663316	13.86%	1.589%
17	7.158	688	692	697	rVB5	6849	10346	0.22%	0.025%
18	7.258	703	709	719	rVB	533023	870701	18.19%	2.086%
19	7.334	719	722	724	rBV2	8938	12403	0.26%	0.030%
20	7.728	782	789	795	rBV	558721	893117	18.66%	2.140%
21	7.793	795	800	808	rVB2	12370	19924	0.42%	0.048%
22	7.963	824	829	836	rBV	12678	20307	0.42%	0.049%
23	8.269	876	881	895	rVB2	10593	21038	0.44%	0.050%
24	8.416	899	906	914	rBV	628121	991007	20.71%	2.375%
25	8.540	921	927	932	rVB2	18080	28988	0.61%	0.069%
26	8.816	969	974	977	rBV2	13285	18794	0.39%	0.045%
27	8.869	977	983	991	rVB	475145	768946	16.07%	1.842%
28	9.587	1094	1105	1112	rBV	349477	572300	11.96%	1.371%
29	10.116	1187	1195	1209	rVB	606423	993800	20.76%	2.381%
30	10.275	1215	1222	1236	rBV	781295	1213966	25.36%	2.909%
31	10.499	1252	1260	1266	rBV	777656	1270468	26.54%	3.044%
32	10.546	1266	1268	1273	rVB	11649	14854	0.31%	0.036%
33	10.628	1275	1282	1293	rBV	561182	938594	19.61%	2.249%
34	11.275	1385	1392	1398	rBV8	4167	8774	0.18%	0.021%
35	12.087	1525	1530	1534	rVB	10054	13934	0.29%	0.033%
36	12.157	1539	1542	1547	rVB	9455	12975	0.27%	0.031%

LSC Area Percent Report

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Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDM5

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_N\METHODS\SFAM-EPA-BN061821MA.M
Title : SVOA CALIBRATION

37	12.381	1575	1580	1585	rBV3	7127	11956	0.25%	0.029%
38	12.704	1629	1635	1639	rBV5	6065	10521	0.22%	0.025%
39	12.898	1664	1668	1674	rVV5	5778	9654	0.20%	0.023%
40	12.957	1674	1678	1687	rVB3	11154	17381	0.36%	0.042%
41	13.187	1710	1717	1722	rBV2	22454	35461	0.74%	0.085%
42	13.251	1725	1728	1733	rVB4	5858	9657	0.20%	0.023%
43	13.551	1776	1779	1784	rVV4	13848	18956	0.40%	0.045%
44	13.669	1795	1799	1803	rVB3	10624	13856	0.29%	0.033%
45	13.757	1808	1814	1822	rVV	1001602	1393909	29.12%	3.340%
46	13.845	1825	1829	1833	rVV4	9649	15520	0.32%	0.037%
47	13.898	1833	1838	1844	rVV7	6210	11360	0.24%	0.027%
48	14.040	1853	1862	1873	rBV	1222362	1827175	38.18%	4.378%
49	14.269	1895	1901	1905	rBV	27650	40759	0.85%	0.098%
50	14.351	1907	1915	1929	rVV	1122283	1659569	34.67%	3.977%
51	14.540	1938	1947	1963	rVV	359907	542651	11.34%	1.300%
52	14.692	1967	1973	1979	rBV7	8633	18635	0.39%	0.045%
53	14.769	1979	1986	1991	rVV3	55523	106147	2.22%	0.254%
54	14.887	2002	2006	2012	rVB7	7200	12533	0.26%	0.030%
55	15.175	2047	2055	2059	rBV7	8638	19017	0.40%	0.046%
56	15.222	2059	2063	2067	rVV	21509	29896	0.62%	0.072%
57	15.257	2067	2069	2075	rVV3	9503	14647	0.31%	0.035%
58	15.345	2075	2084	2090	rVV	1526811	2182271	45.60%	5.229%
59	15.398	2090	2093	2098	rVB2	40224	52407	1.09%	0.126%
60	15.457	2098	2103	2109	rBV	258985	360075	7.52%	0.863%
61	15.510	2109	2112	2118	rVB	42167	58935	1.23%	0.141%
62	15.587	2118	2125	2128	rBV	16391	28927	0.60%	0.069%
63	15.698	2141	2144	2146	rBV2	8042	10787	0.23%	0.026%
64	15.775	2151	2157	2165	rBV	198070	278329	5.82%	0.667%
65	15.839	2165	2168	2174	rVB3	15777	27308	0.57%	0.065%
66	15.910	2176	2180	2184	rBV6	5607	12120	0.25%	0.029%
67	16.039	2198	2202	2206	rVB2	17807	23827	0.50%	0.057%
68	16.081	2206	2209	2212	rBV2	22407	27731	0.58%	0.066%
69	16.251	2234	2238	2243	rBV6	8468	14910	0.31%	0.036%
70	16.604	2293	2298	2299	rBV5	9540	15877	0.33%	0.038%
71	16.663	2305	2308	2314	rVB	22894	34091	0.71%	0.082%
72	16.845	2334	2339	2341	rBV6	12675	22394	0.47%	0.054%
73	16.875	2341	2344	2348	rVV2	38545	49593	1.04%	0.119%
74	16.922	2348	2352	2357	rVB3	22899	42464	0.89%	0.102%
75	17.098	2376	2382	2386	rBV	1322857	1925067	40.22%	4.613%
76	17.139	2386	2389	2393	rVV	226393	301298	6.30%	0.722%

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Acq On : 26 Jun 2021 15:12
Operator : CG/JU
Sample : M2704-15
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDM5

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

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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Title : SVOA CALIBRATION

77	17.198	2393	2399	2411	rVV	1575857	2363856	49.39%	5.664%
78	17.292	2412	2415	2421	rVB6	28638	44074	0.92%	0.106%
79	17.351	2421	2425	2427	rBV4	11844	18015	0.38%	0.043%
80	17.475	2442	2446	2448	rBV4	14954	23243	0.49%	0.056%
81	17.616	2466	2470	2477	rBV	45125	61873	1.29%	0.148%
82	17.786	2494	2499	2504	rBV	157429	196751	4.11%	0.471%
83	17.975	2528	2531	2532	rBV	21695	21909	0.46%	0.052%
84	18.004	2532	2536	2542	rVB3	63588	120446	2.52%	0.289%
85	18.251	2575	2578	2581	rBV2	18784	19523	0.41%	0.047%
86	18.310	2584	2588	2597	rBV2	64111	101549	2.12%	0.243%
87	18.422	2603	2607	2612	rBV	173513	244527	5.11%	0.586%
88	18.963	2692	2699	2705	rVB2	94878	188139	3.93%	0.451%
89	19.163	2729	2733	2739	rVB	209969	282393	5.90%	0.677%
90	19.498	2784	2790	2794	rBV	1888575	2764250	57.76%	6.623%
91	20.063	2883	2886	2890	rVB	49207	53464	1.12%	0.128%
92	20.516	2959	2963	2968	rBV	4035348	4786108	100.00%	11.468%
93	21.027	3047	3050	3055	rBV3	135044	189949	3.97%	0.455%
94	21.222	3080	3083	3089	rVB	257016	305372	6.38%	0.732%
95	21.298	3090	3096	3100	rBV	1634276	2191470	45.79%	5.251%
96	22.163	3239	3243	3247	rBV	182806	240803	5.03%	0.577%
97	22.798	3348	3351	3357	rVB2	90366	128580	2.69%	0.308%
98	23.404	3447	3454	3460	rBV2	1292968	2436023	50.90%	5.837%
99	23.545	3471	3478	3492	rBV	1153591	2147504	44.87%	5.146%
100	24.104	3570	3573	3580	rVB	105844	180295	3.77%	0.432%

Sum of corrected areas: 41734118

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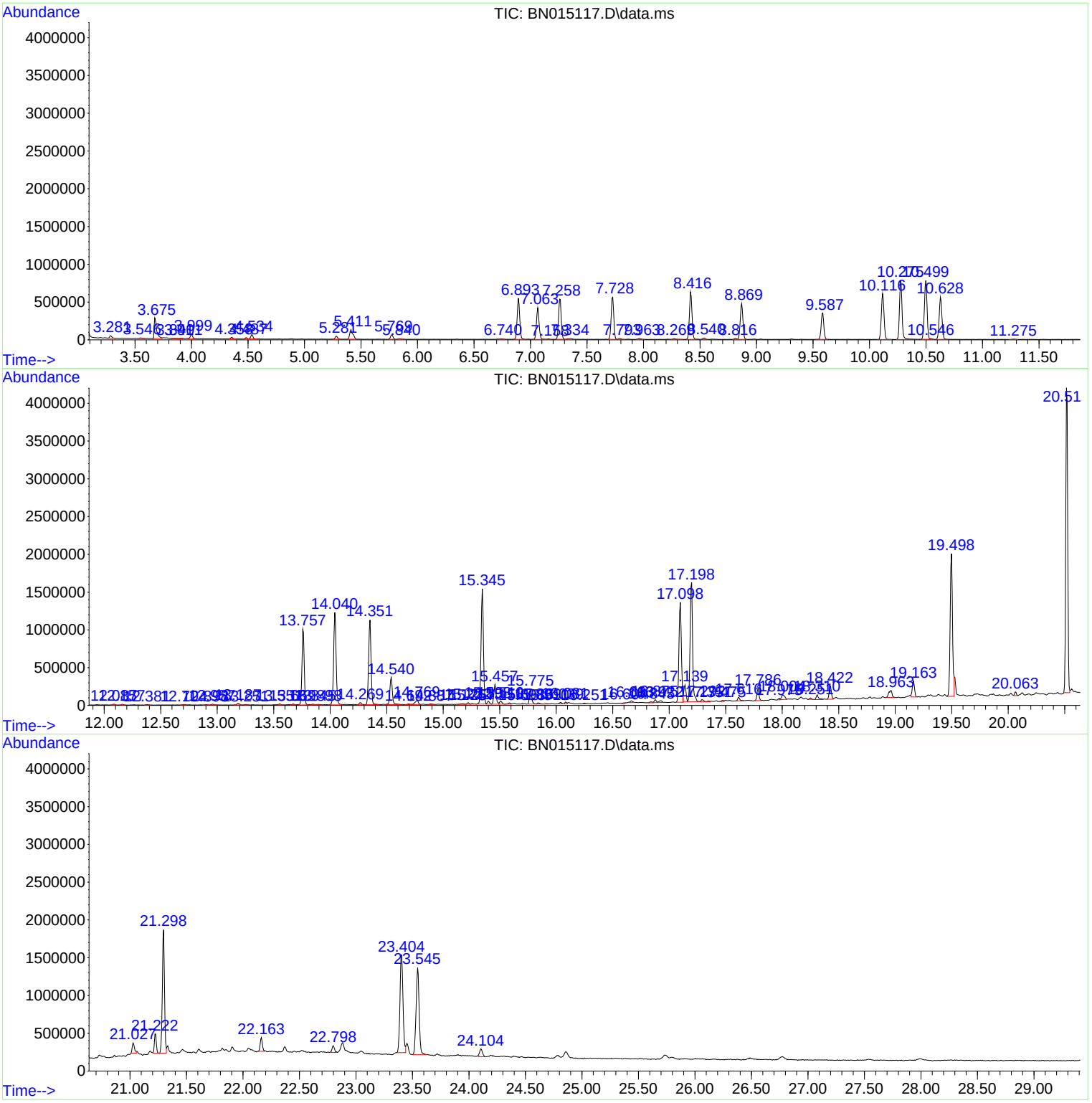
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDM5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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Data File : BN015117.D
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Instrument :
BNA_N
ClientSampleId :
BGDM5

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Quant Title : SVOA CALIBRATION

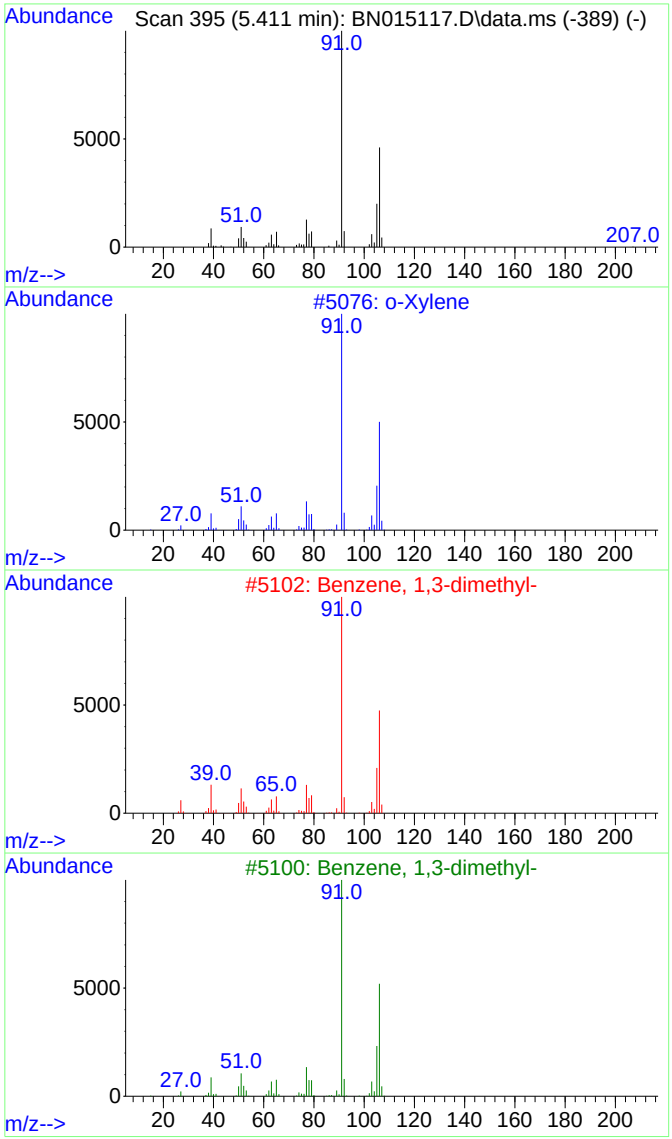
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

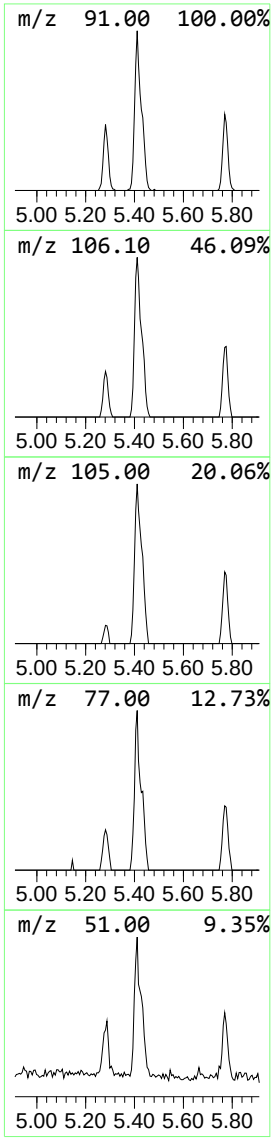
****

Peak Number 1 o-Xylene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.411	5.46 ng/ul	243814	1,4-Dichlorobenzene-d4	7.728

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Xylene	106	C8H10	000095-47-6	97
2	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
3	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
4	p-Xylene	106	C8H10	000106-42-3	97
5	p-Xylene	106	C8H10	000106-42-3	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015117.D
Acq On : 26 Jun 2021 15:12
Operator : CG/JU
Sample : M2704-15
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDM5

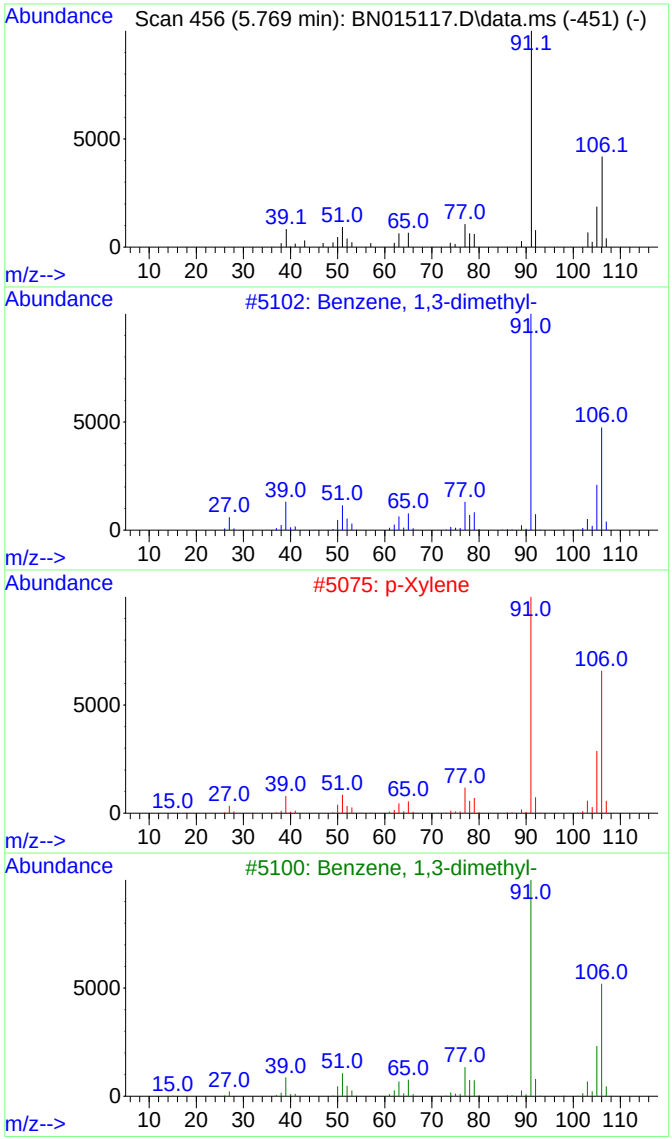
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

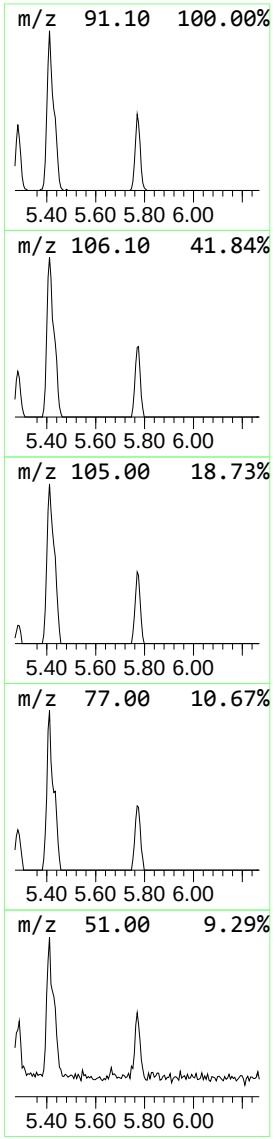
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.769	2.02 ng/ul	90007	1,4-Dichlorobenzene-d4	7.728

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
2	p-Xylene	106	C8H10	000106-42-3	95
3	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94
4	o-Xylene	106	C8H10	000095-47-6	94
5	p-Xylene	106	C8H10	000106-42-3	94





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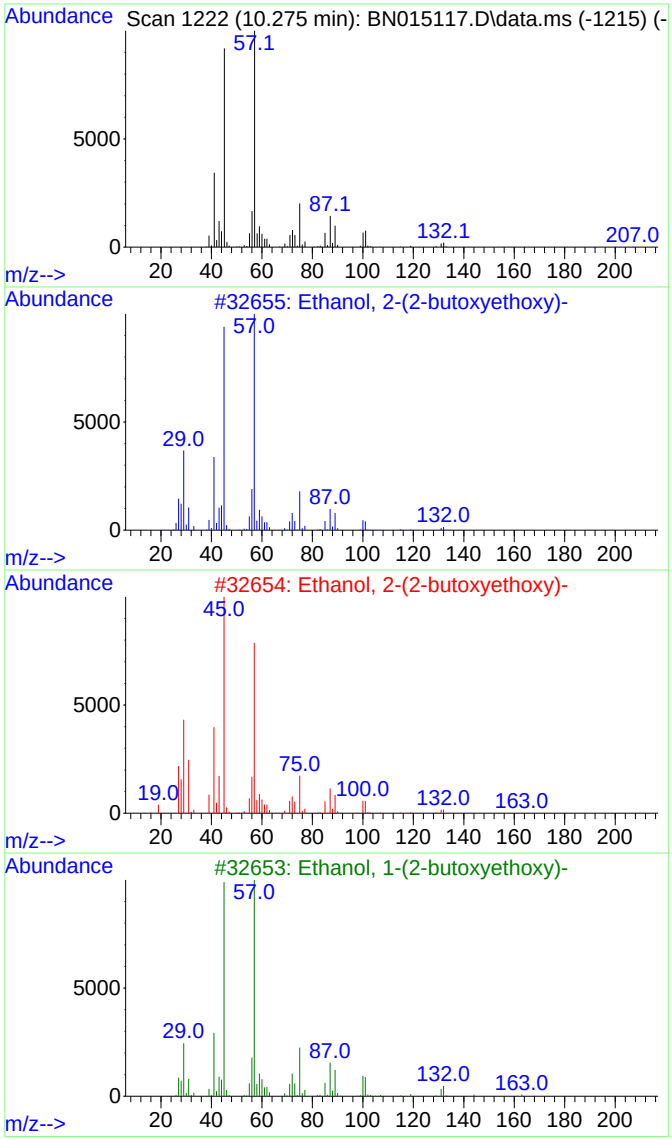
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

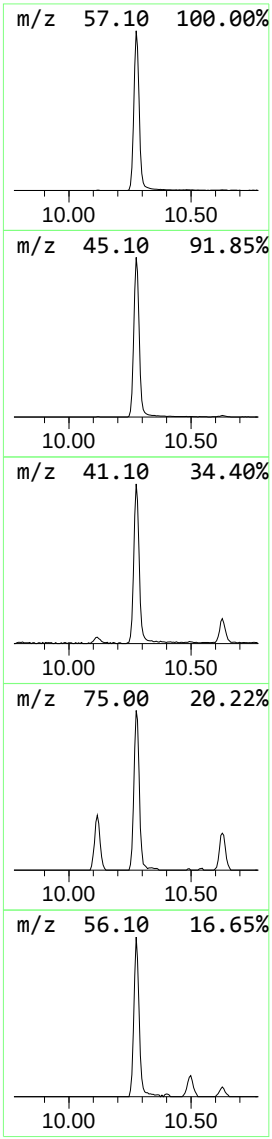
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.275	19.11 ng/ul	1213970	Naphthalene-d8	10.499

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72





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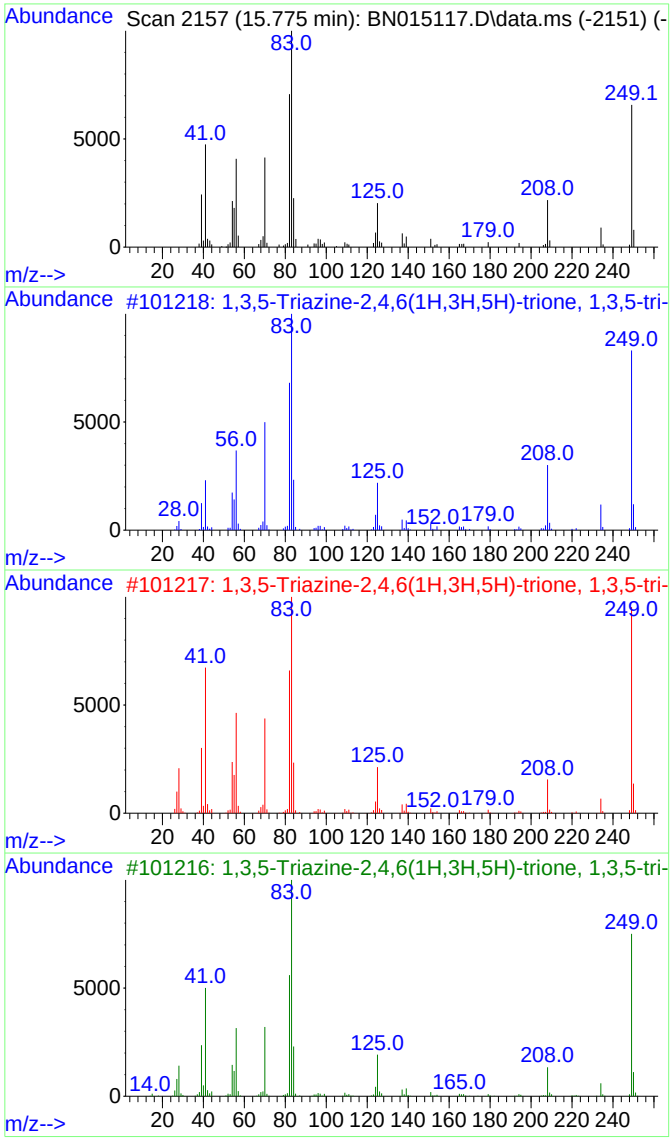
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

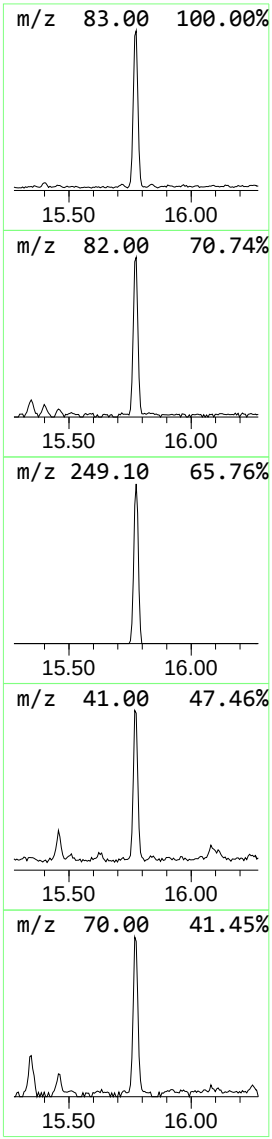
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1,3,5-Triazine-2,4,6(1H,3H,... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.775	2.89 ng/ul	278329	Phenanthrene-d10	17.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	98
2	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	96
3	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	95
4	N-Carboxy-1-[4-chlorophenyl]-2-[...	295	C15H18ClNO3	058158-38-6	37
5	Hexane, 1,6-dicyclohexyl-	250	C18H34	001610-23-7	27





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Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDM5

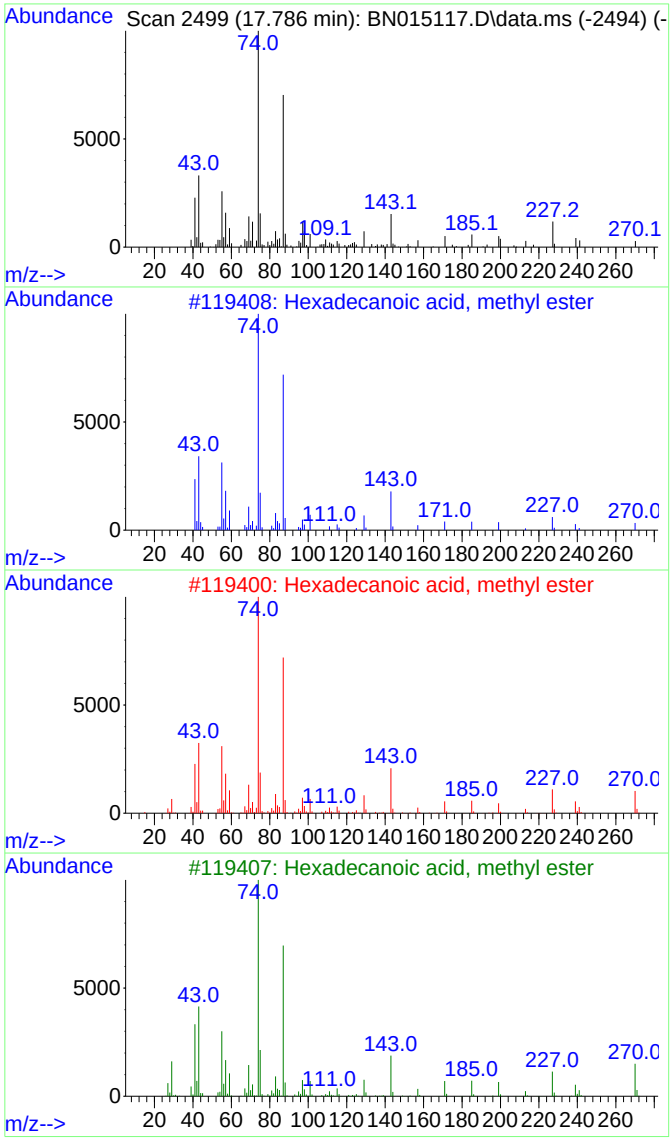
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

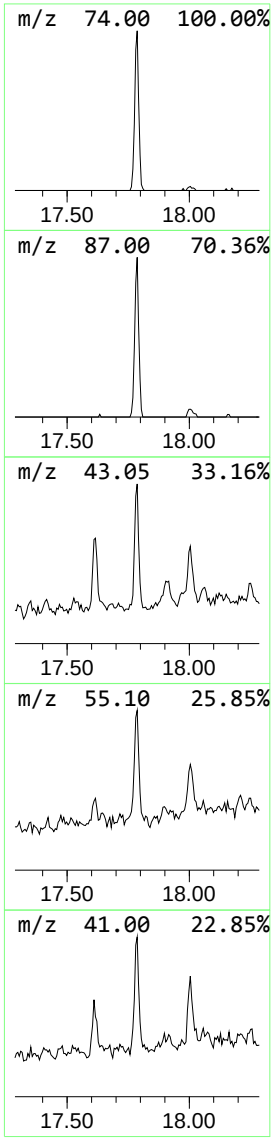
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.786	2.04 ng/ul	196751	Phenanthrene-d10	17.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
3	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
4	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	81
5	Undecanoic acid, methyl ester	200	C12H24O2	001731-86-8	81





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Acq On : 26 Jun 2021 15:12
Operator : CG/JU
Sample : M2704-15
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDM5

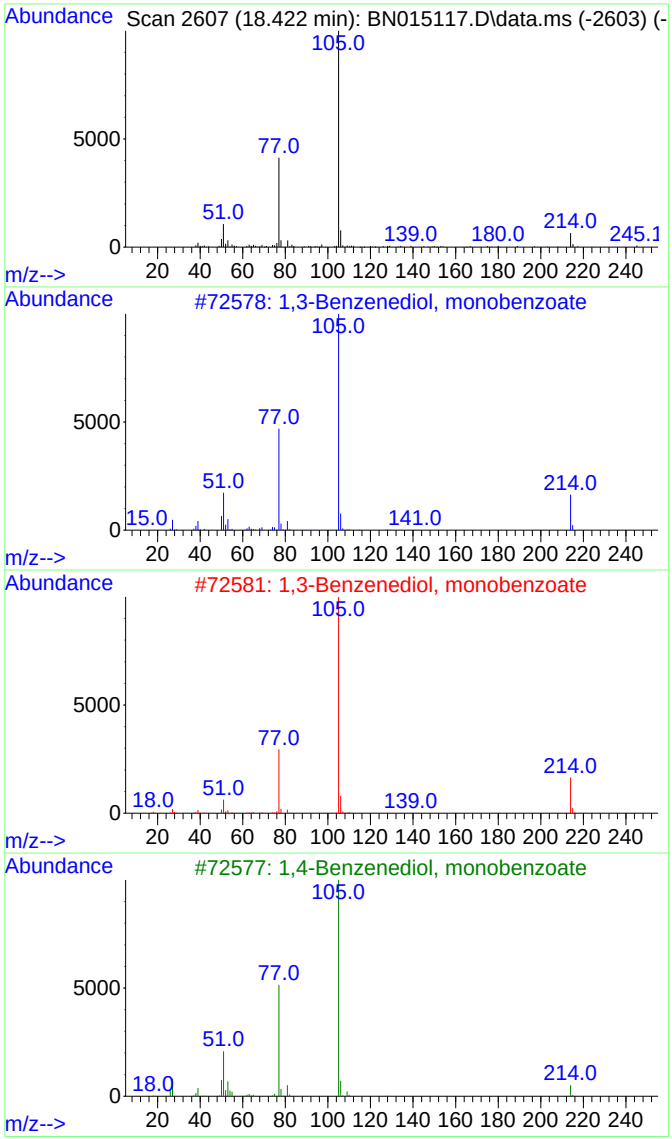
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

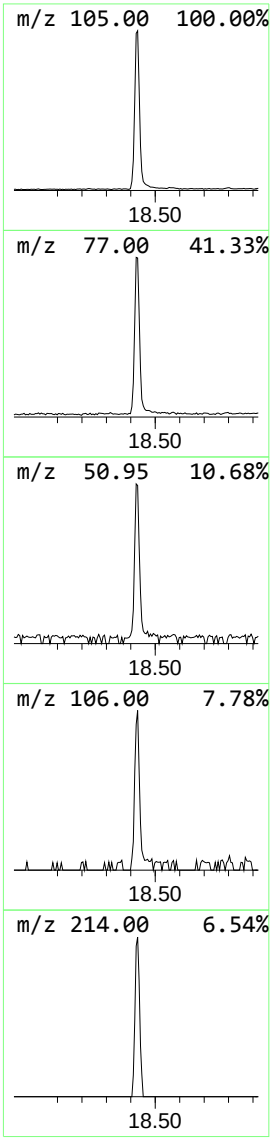
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 1,3-Benzenediol, monobenzoate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.422	2.54 ng/ul	244527	Phenanthrene-d10	17.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	90
2	1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	90
3	1,4-Benzenediol, monobenzoate	214	C13H10O3	002444-19-1	87
4	1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	86
5	Isopropyl phenyl ketone	148	C10H12O	000611-70-1	72





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Data File : BN015117.D
Acq On : 26 Jun 2021 15:12
Operator : CG/JU
Sample : M2704-15
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDM5

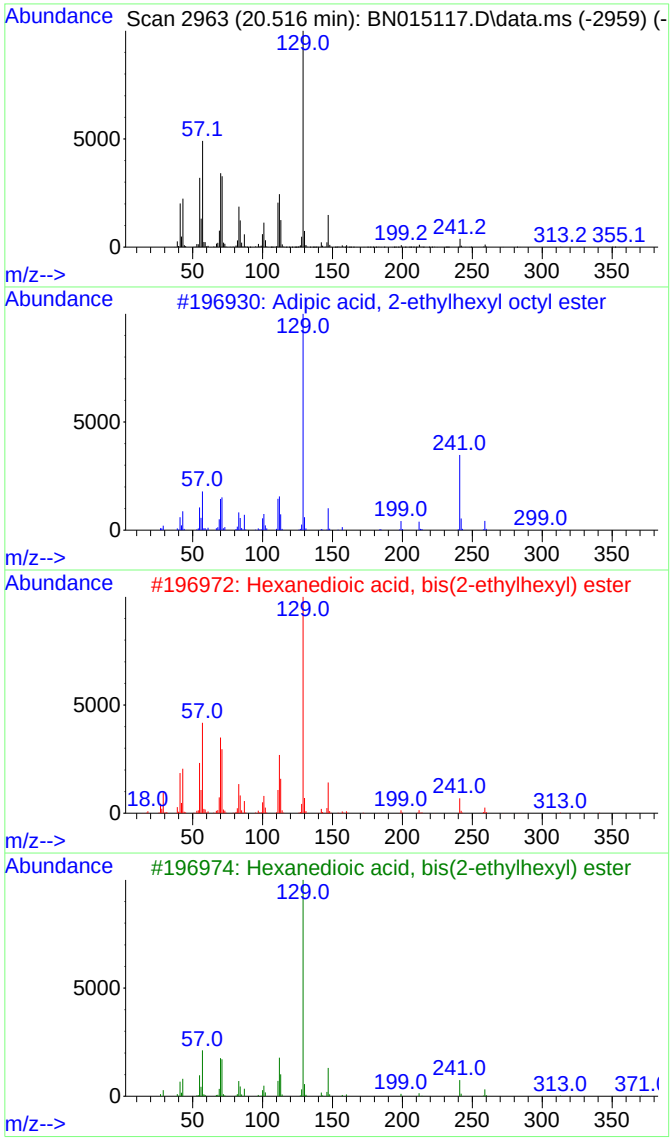
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

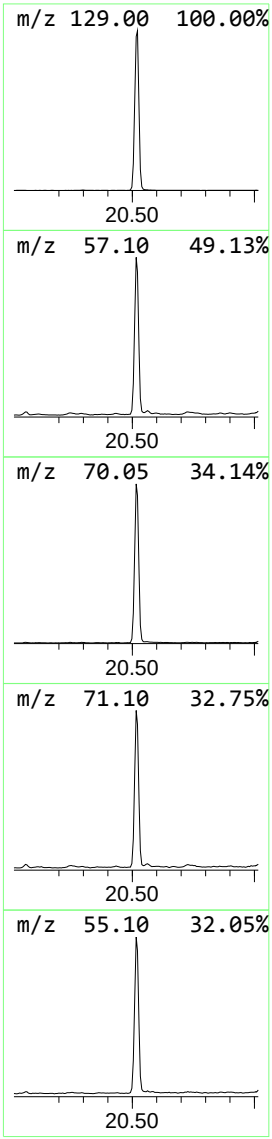
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Adipic acid, 2-ethylhexyl o... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.516	43.68 ng/ul	4786110	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Adipic acid, 2-ethylhexyl octyl ...	370	C22H42O4	1000324-48-6	91
2	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	90
3	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
4	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	86
5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	76





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015117.D
Acq On : 26 Jun 2021 15:12
Operator : CG/JU
Sample : M2704-15
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDM5

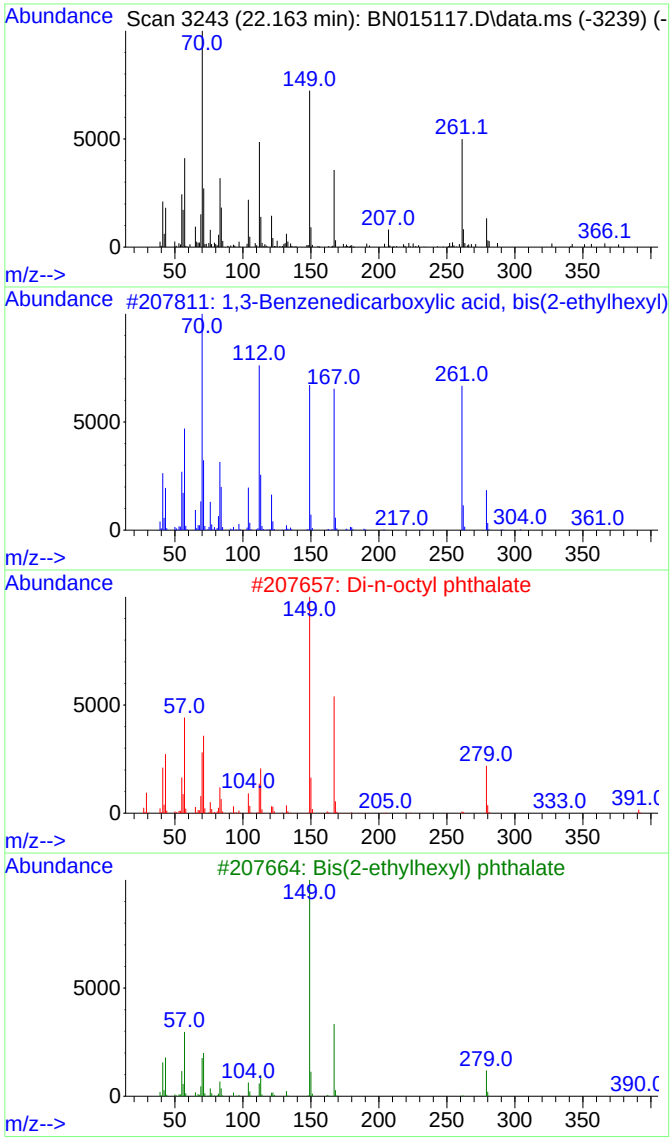
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

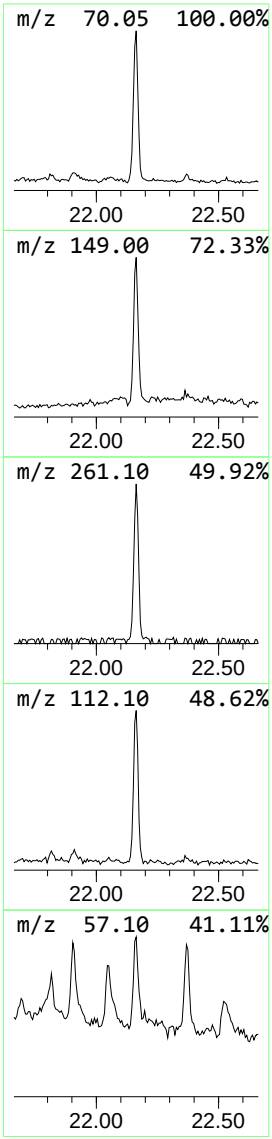
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1,3-Benzenedicarboxylic aci... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.163	2.20 ng/ul	240803	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	62
2	Di-n-octyl phthalate	390	C24H38O4	000117-84-0	25
3	Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	22
4	Phthalic acid, 3-methoxybenzyl p...	356	C21H24O5	1000377-97-8	11
5	Diethyl Phthalate	222	C12H14O4	000084-66-2	11





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015117.D
Acq On : 26 Jun 2021 15:12
Operator : CG/JU
Sample : M2704-15
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDM5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
o-Xylene	5.411	5.5	ng/ul	243814	1	7.728	893117	20.0
Benzene, 1,3-di...	5.769	2.0	ng/ul	90007	1	7.728	893117	20.0
Ethanol, 2-(2-b...	10.275	19.1	ng/ul	1213970	2	10.499	1270470	20.0
1,3,5-Triazine-...	15.775	2.9	ng/ul	278329	4	17.098	1925070	20.0
Hexadecanoic ac...	17.786	2.0	ng/ul	196751	4	17.098	1925070	20.0
1,3-Benzenediol...	18.422	2.5	ng/ul	244527	4	17.098	1925070	20.0
Adipic acid, 2-...	20.516	43.7	ng/ul	4786110	5	21.298	2191470	20.0
1,3-Benzenedica...	22.163	2.2	ng/ul	240803	5	21.298	2191470	20.0