LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015103.D
Acq On : 26 Jun 2021 06:51
Operator : CG/JU
Sample : M2704-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDB7

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: BN015103.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	44	rVB	34849	50835	0.47%	0.107%
2	3.675	97	100	112	rBV	316163	466127	4.27%	0.979%
3	3.999	151	155	162	rVB	56828	80960	0.74%	0.170%
4	4.358	211	216	220	rVB	19389	26721	0.24%	0.056%
5	4.481	233	237	242	rBV	16367	22646	0.21%	0.048%
6	4.534	242	246	250	rVB	50522	64658	0.59%	0.136%
7	5.281	368	373	380	rBV	38893	54646	0.50%	0.115%
8	5.411	389	395	405	rVB	117769	236769	2.17%	0.497%
9	5.769	451	456	463	rBV	55455	84360	0.77%	0.177%
10	5.840	464	468	473	rVB2	8253	11994	0.11%	0.025%
11	6.893	640	647	656	rBV	518767	802604	7.35%	1.686%
12	7.063	668	676	685	rBV	400281	629549	5.76%	1.322%
13	7.163	689	693	697	rVB3	7200	10939	0.10%	0.023%
14	7.258	702	709	722	rVB	521990	836402	7.66%	1.757%
15	7.358	722	726	731	rBV4	6749	11429	0.10%	0.024%
16	7.728	782	789	795	rBV	638113	991043	9.07%	2.082%
17	7.793	795	800	805	rVB	18123	28441	0.26%	0.060%
18	7.963	825	829	837	rVB3	10135	18698	0.17%	0.039%
19	8.416	898	906	915	rBV	589095	928674	8.50%	1.951%
20	8.540	921	927	931	rVB	16499	27061	0.25%	0.057%
21	8.810	968	973	977	rBV2	14069	23746	0.22%	0.050%
22	8.869	977	983	991	rVB	450117	733596	6.71%	1.541%
23	9.587	1098	1105	1112	rVB	322584	524797	4.80%	1.102%
24	10.116	1189	1195	1205	rBV	577810	938244	8.59%	1.971%
25	10.275	1217	1222	1231	rBV	135879	220171	2.02%	0.462%
26	10.499	1252	1260	1267	rBV	877226	1418127	12.98%	2.979%
27	10.628	1275	1282	1290	rBV	557228	920824	8.43%	1.934%
28	12.087	1521	1530	1536	rVB	9454	17127	0.16%	0.036%
29	13.187	1713	1717	1723	rVB2	16901	23351	0.21%	0.049%
30	13.763	1808	1815	1820	rBV	936672	1314129	12.03%	2.760%
31	14.040	1852	1862	1869	rBV	1169467	1686530	15.44%	3.543%
32	14.269	1897	1901	1905	rVB	22966	27144	0.25%	0.057%
33	14.351	1909	1915	1930	rVB	1277117	1830994	16.76%	3.846%
34	14.540	1939	1947	1957	rBV	315258	466173	4.27%	0.979%
35	14.769	1981	1986	1990	rVB	40747	54335	0.50%	0.114%
36	15.222	2059	2063	2067	rBV	14575	18303	0.17%	0.038%

LSC Area Percent Report

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Instrument : BNA_N
ClientSampleId : BGDB7

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	15.345	2078	2084	2091	rVB	1454351	2060021	18.86%	4.327%
38	15.457	2098	2103	2109	rBV	212404	297376	2.72%	0.625%
39	15.510	2109	2112	2116	rVB	36606	46936	0.43%	0.099%
40	15.587	2121	2125	2130	rBV2	14508	21352	0.20%	0.045%
41	15.775	2151	2157	2164	rBV	199713	258075	2.36%	0.542%
42	16.081	2206	2209	2213	rBV2	9344	12329	0.11%	0.026%
43	16.251	2233	2238	2243	rBV5	10256	14571	0.13%	0.031%
44	16.604	2293	2298	2301	rBV6	6829	11369	0.10%	0.024%
45	16.651	2301	2306	2310	rBV6	5905	11075	0.10%	0.023%
46	16.875	2340	2344	2348	rVB3	24055	33838	0.31%	0.071%
47	17.098	2376	2382	2388	rBV	1518315	2122670	19.43%	4.459%
48	17.192	2392	2398	2407	rVV2	1519679	2255258	20.64%	4.737%
49	17.292	2411	2415	2421	rVB5	21952	33957	0.31%	0.071%
50	17.345	2421	2424	2427	rBV5	7995	11157	0.10%	0.023%
51	17.616	2466	2470	2474	rBV	30898	44576	0.41%	0.094%
52	17.786	2496	2499	2506	rVB6	29689	47610	0.44%	0.100%
53	18.004	2532	2536	2542	rBV5	31697	46462	0.43%	0.098%
54	18.245	2575	2577	2582	rVB2	19723	23456	0.21%	0.049%
55	18.310	2584	2588	2596	rBV	41821	60703	0.56%	0.128%
56	18.428	2603	2608	2623	rVB	350450	500635	4.58%	1.052%
57	18.728	2655	2659	2663	rVB3	27889	39476	0.36%	0.083%
58	18.945	2691	2696	2698	rBV2	73143	109261	1.00%	0.230%
59	19.128	2724	2727	2731	rVB	22446	24304	0.22%	0.051%
60	19.210	2739	2741	2745	rBV5	19486	22164	0.20%	0.047%
61	19.286	2751	2754	2758	rBV5	20615	31933	0.29%	0.067%
62	19.380	2767	2770	2774	rVB3	16987	21788	0.20%	0.046%
63	19.498	2784	2790	2799	rBV	1935109	2800826	25.64%	5.883%
64	20.063	2882	2886	2889	rBV2	37827	45232	0.41%	0.095%
65	20.198	2906	2909	2914	rBV6	29348	45279	0.41%	0.095%
66	20.245	2914	2917	2922	rVV6	20244	38528	0.35%	0.081%
67	20.516	2958	2963	2969	rBV	8983615	10924971	100.00%	22.948%
68	20.857	3019	3021	3025	rBV5	23066	24998	0.23%	0.053%
69	21.022	3046	3049	3052	rBV	67325	82528	0.76%	0.173%
70	21.174	3072	3075	3078	rBV3	37374	51530	0.47%	0.108%
71	21.222	3079	3083	3089	rVB	287930	329341	3.01%	0.692%
72	21.292	3089	3095	3100	rBV	1847912	2371891	21.71%	4.982%
73	21.904	3195	3199	3209	rVB2	77727	132209	1.21%	0.278%
74	22.157	3238	3242	3246	rVB	195193	244249	2.24%	0.513%
75	22.363	3274	3277	3282	rVB	80165	97278	0.89%	0.204%
76	22.792	3346	3350	3358	rVB	205758	320095	2.93%	0.672%

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Instrument : BNA_N
ClientSampleId : BGDB7

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
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Title : SVOA CALIBRATION

77	22.874	3360	3364	3370	rVB	151383	212248	1.94%	0.446%
78	23.392	3445	3452	3458	rBV	1464415	2734826	25.03%	5.745%
79	23.445	3458	3461	3467	rVV2	157325	272089	2.49%	0.572%
80	23.539	3470	3477	3488	rVB2	1306761	2503514	22.92%	5.259%
81	24.098	3567	3572	3579	rVB2	159004	298915	2.74%	0.628%
82	24.851	3695	3700	3708	rVB	139412	311938	2.86%	0.655%

Sum of corrected areas: 47606984

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ALS Vial : 32 Sample Multiplier: 1

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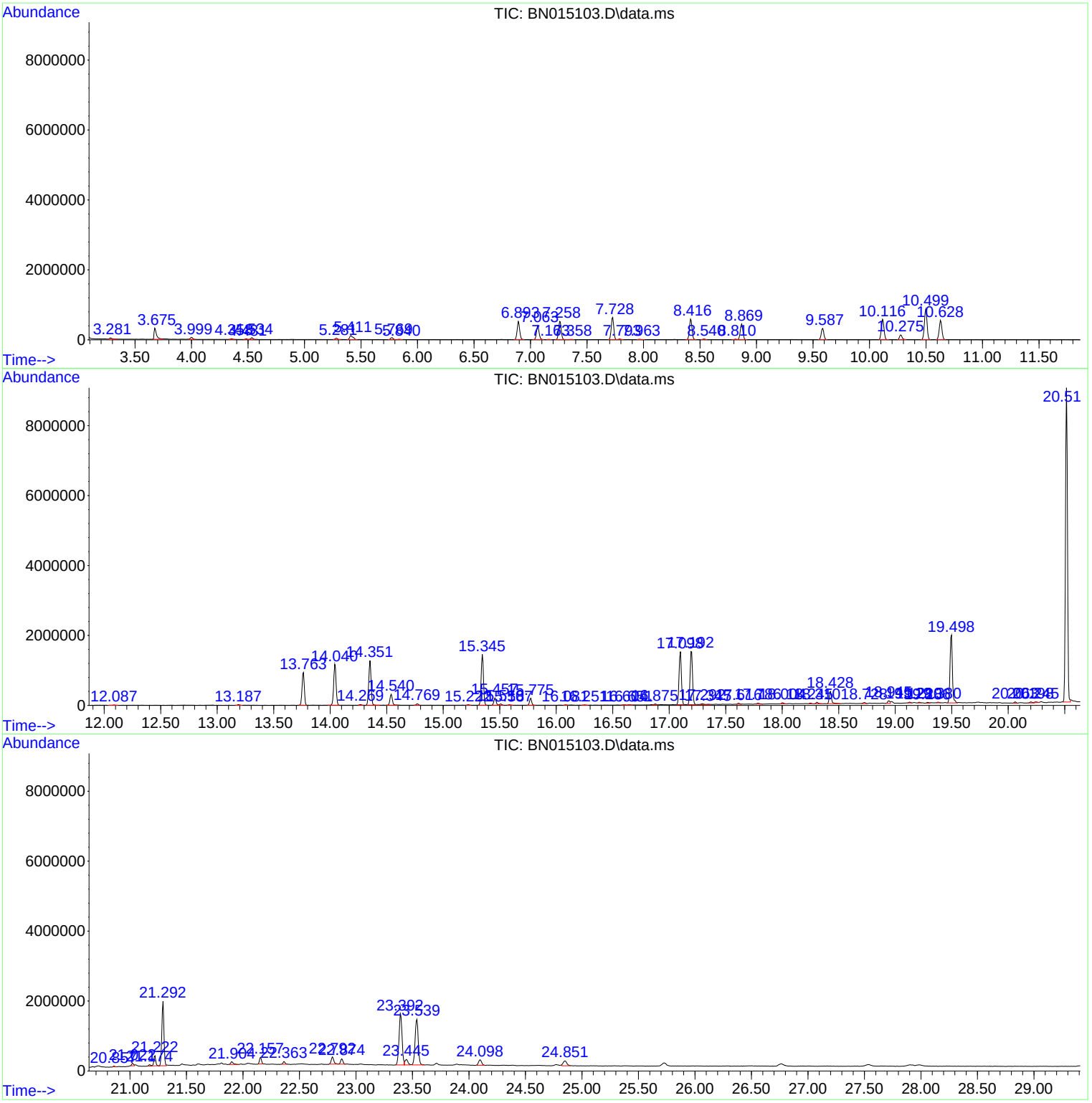
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015103.D
Acq On : 26 Jun 2021 06:51
Operator : CG/JU
Sample : M2704-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDB7

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

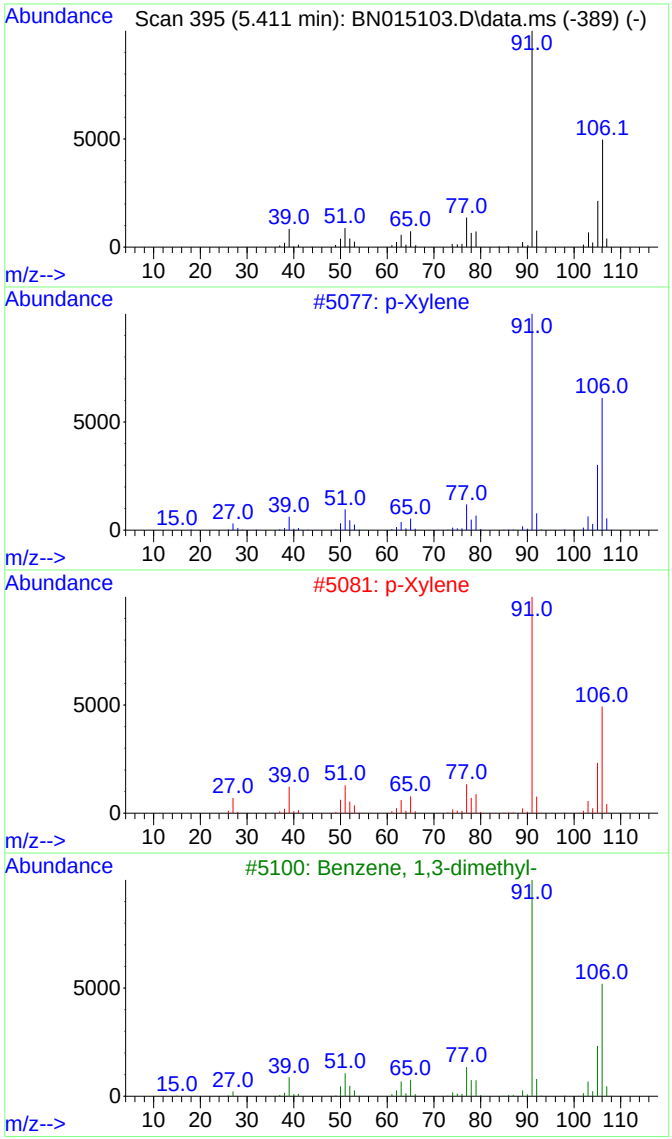
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

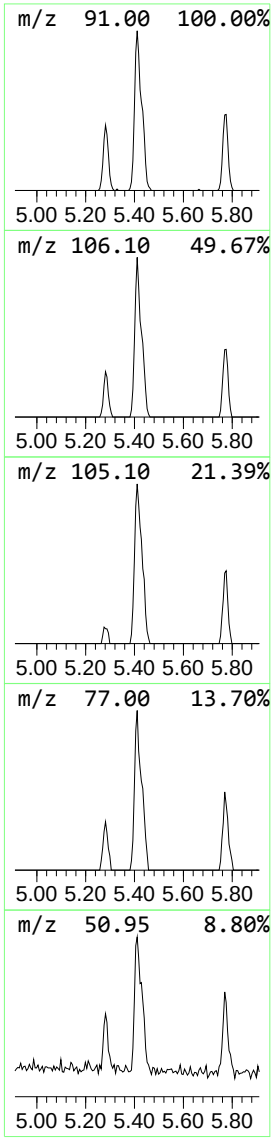
****

Peak Number 1 p-Xylene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.411	4.78 ng/ul	236769	1,4-Dichlorobenzene-d4	7.728

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





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
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Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDB7

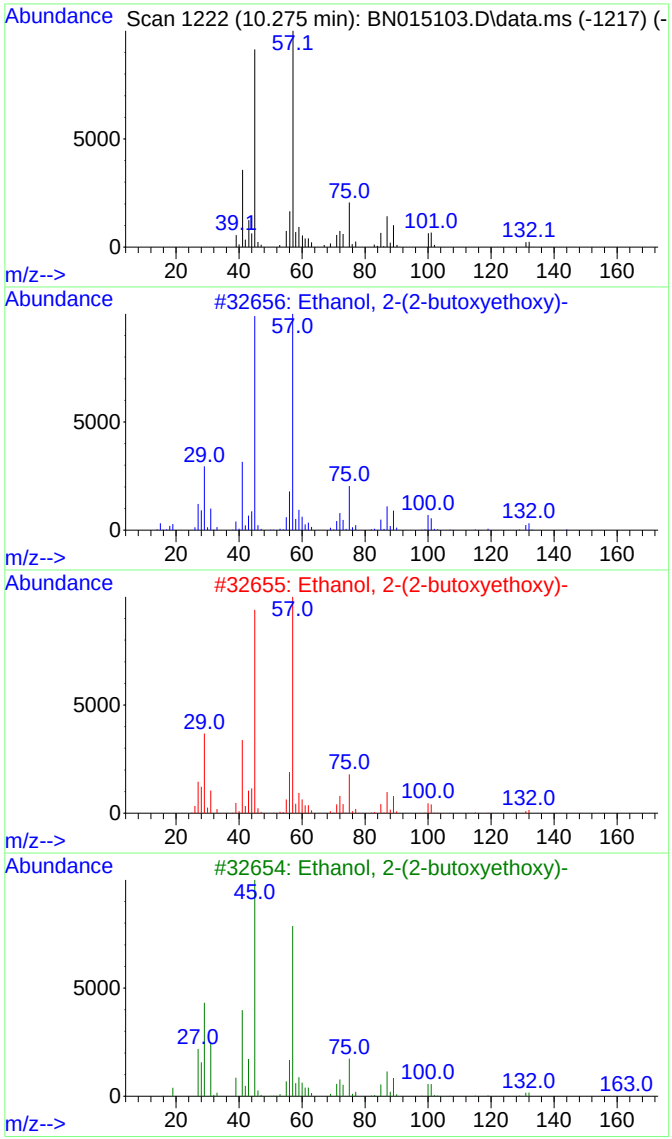
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

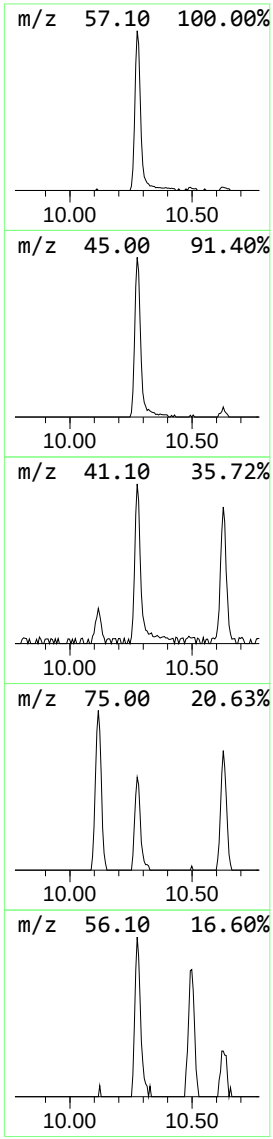
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.275	3.11 ng/ul	220171	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, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
5	Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
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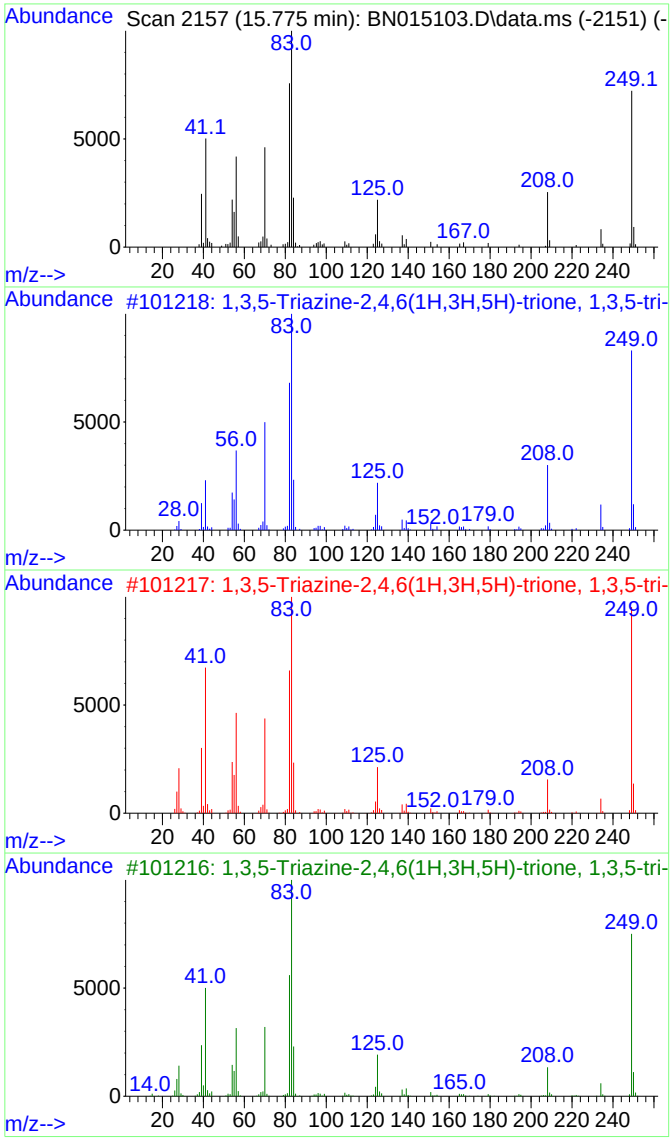
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

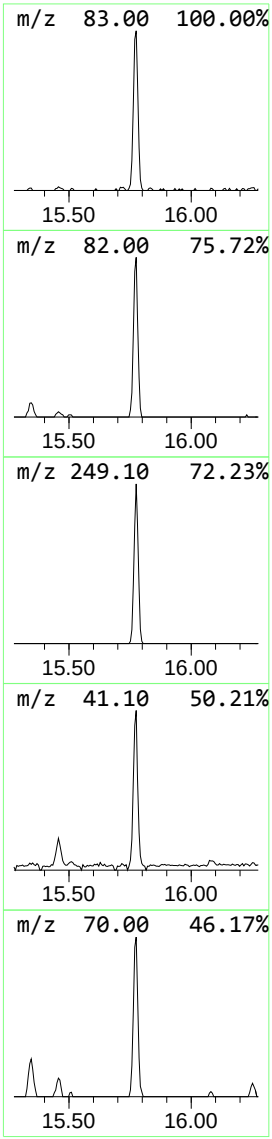
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1,3,5-Triazine-2,4,6(1H,3H,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.775	2.43 ng/ul	258075	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	99
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	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	25
5	Triallyl cyanurate	249	C12H15N3O3	000101-37-1	25





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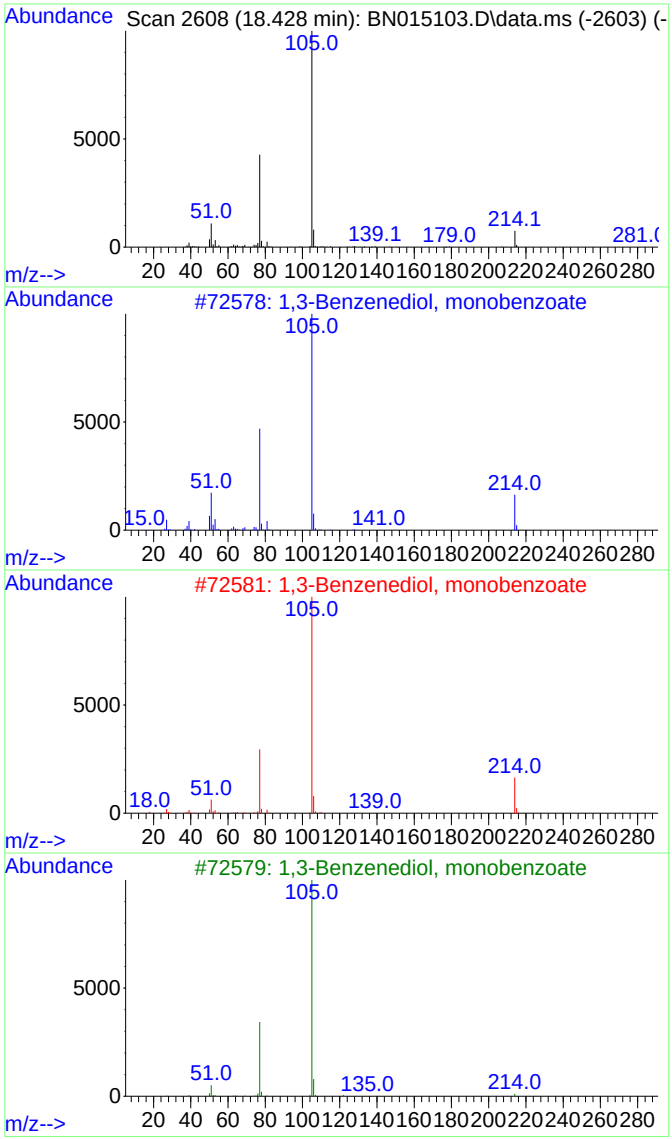
Quant Title : SVOA CALIBRATION

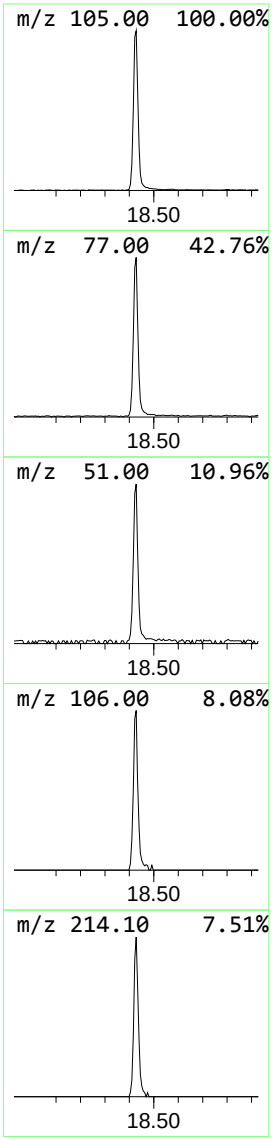
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1,3-Benzenediol, monobenzoate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.427	4.72 ng/ul	500635	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	1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	90
4	2,5-Pyrrolidinedione, 1-(benzoyl...	219	C11H9NO4	023405-15-4	72
5	Benzeneacetic acid, .alpha.-oxo-...	178	C10H10O3	001603-79-8	72





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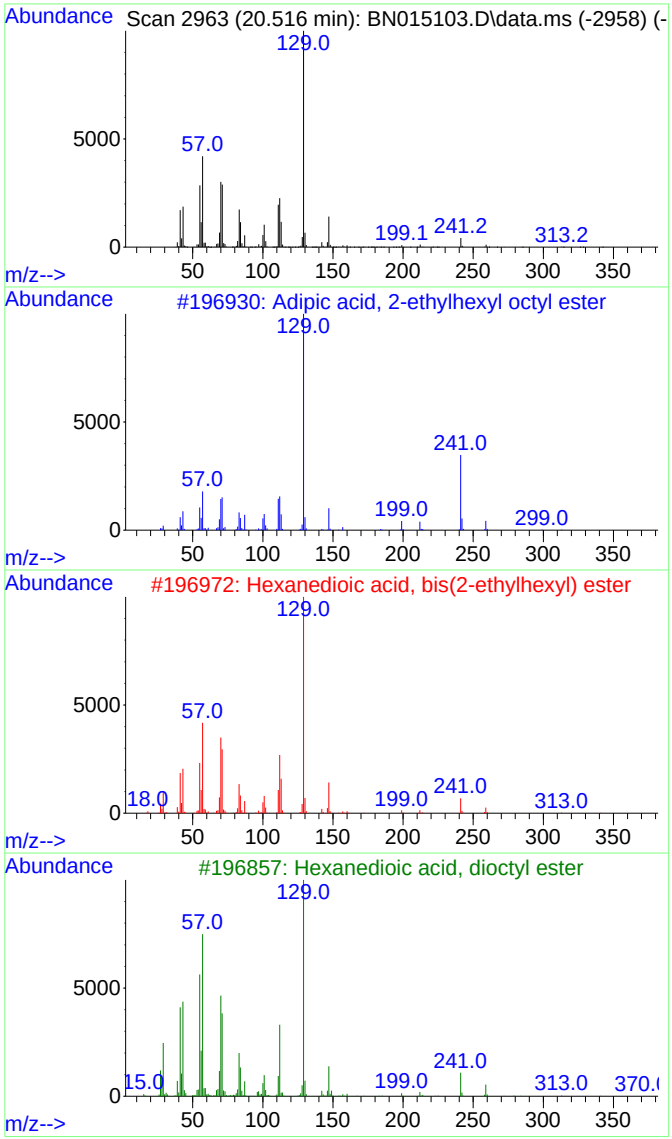
Quant Title : SVOA CALIBRATION

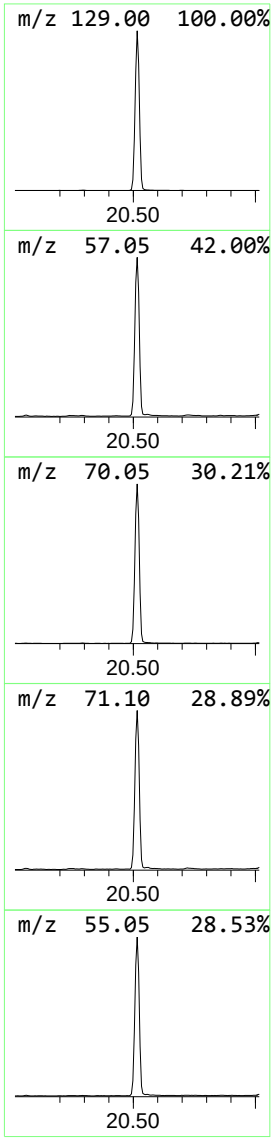
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Adipic acid, 2-ethylhexyl o... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.516	92.12 ng/ul	10925000	Chrysene-d12	21.292

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, dioctyl ester	370	C22H42O4	000123-79-5	87
4	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	86





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ALS Vial : 32 Sample Multiplier: 1

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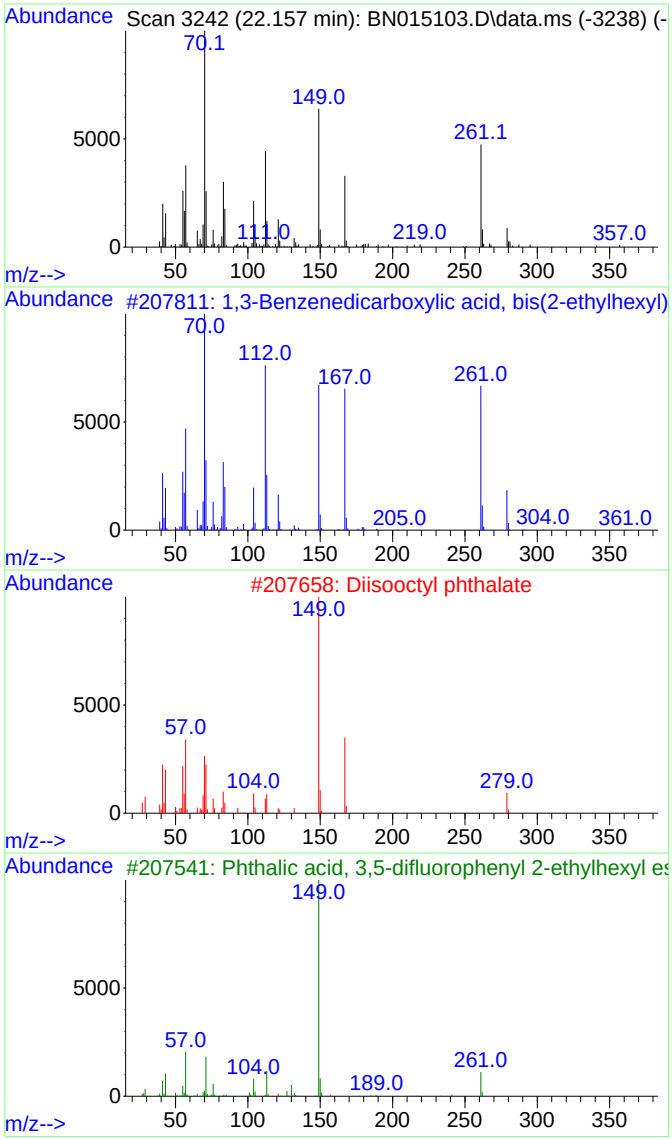
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

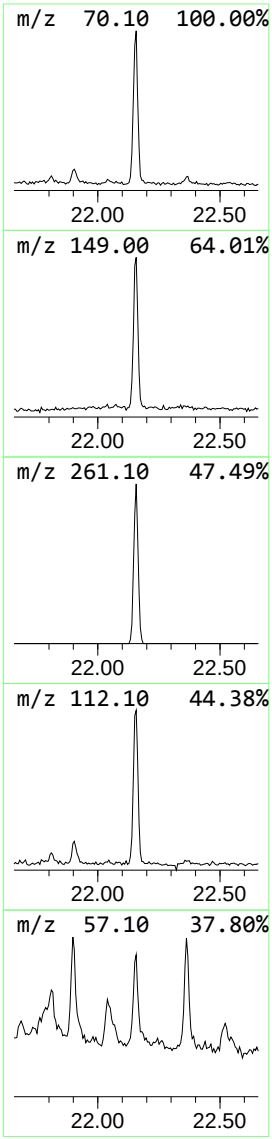
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 1,3-Benzenedicarboxylic aci... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.157	2.06 ng/ul	244249	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	52
2	Diisooctyl phthalate	390	C24H38O4	000131-20-4	22
3	Phthalic acid, 3,5-difluoropheny...	390	C22H24F2O4	1000315-51-3	22
4	Phthalic acid, octyl phenyl ester	354	C22H26O4	1000314-87-4	22
5	Phthalic acid, di(6-methylhept-2...	390	C24H38O4	1000377-97-3	14





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Data File : BN015103.D
Acq On : 26 Jun 2021 06:51
Operator : CG/JU
Sample : M2704-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDB7

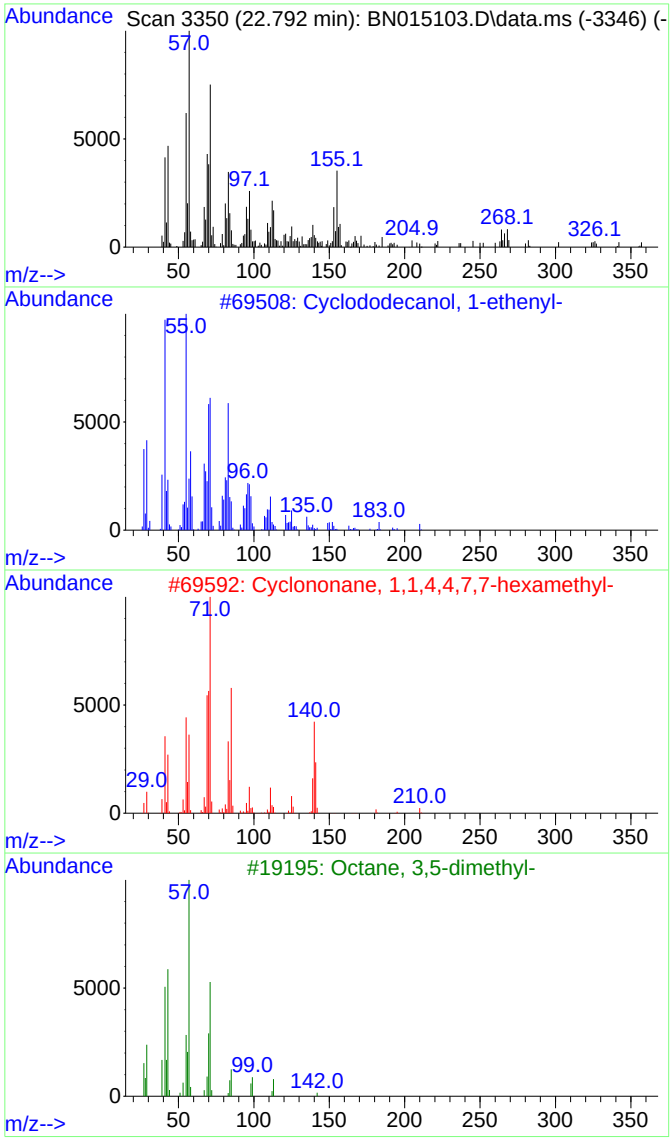
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

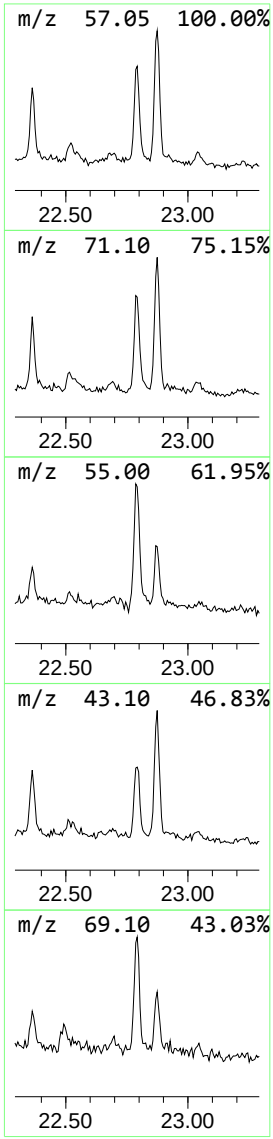
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Cyclododecanol, 1-ethenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.792	2.56 ng/ul	320095	Perylene-d12	23.539

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclododecanol, 1-ethenyl-	210	C14H26O	006244-49-1	91
2	Cyclononane, 1,1,4,4,7,7-hexamet...	210	C15H30	149331-19-1	49
3	Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	46
4	5-(1-Iodo-1-methyl-ethyl)-3,3-di...	282	C9H15IO2	1000190-61-9	43
5	2-Nonen-1-ol, 2-methyl-	156	C10H20O	091008-40-1	30





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Data File : BN015103.D
Acq On : 26 Jun 2021 06:51
Operator : CG/JU
Sample : M2704-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDB7

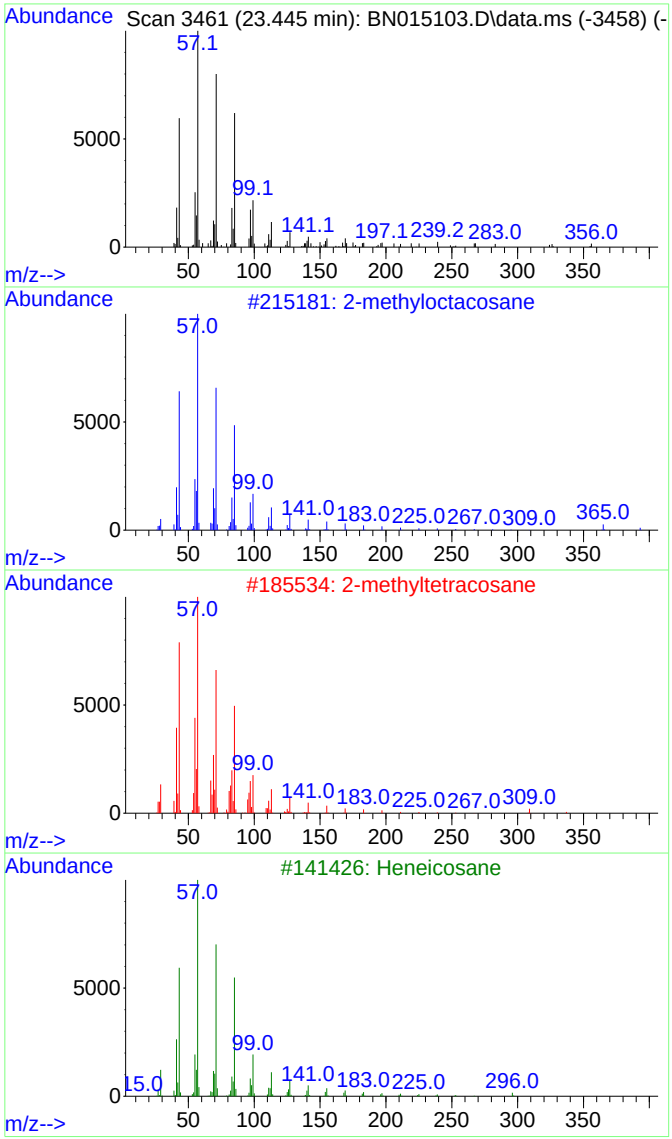
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

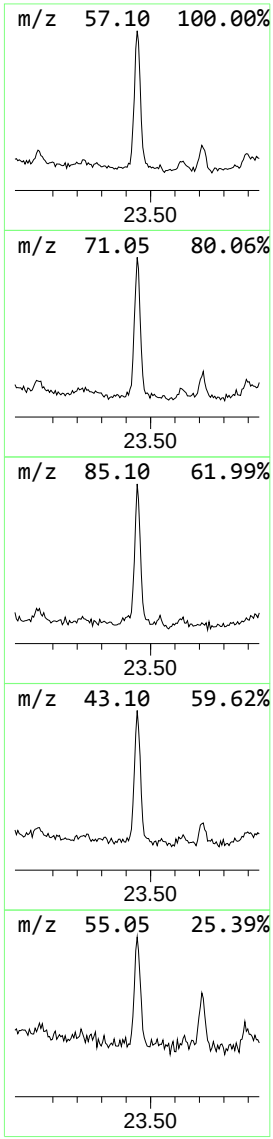
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.445	2.17 ng/ul	272089	Perylene-d12	23.539

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-methyloctacosane	408	C29H60	1000376-72-8	91
2	2-methyltetracosane	352	C25H52	1000376-72-6	91
3	Heneicosane	296	C21H44	000629-94-7	90
4	Triacontane	422	C30H62	000638-68-6	90
5	Hexacosane	366	C26H54	000630-01-3	87





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Data File : BN015103.D
Acq On : 26 Jun 2021 06:51
Operator : CG/JU
Sample : M2704-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDB7

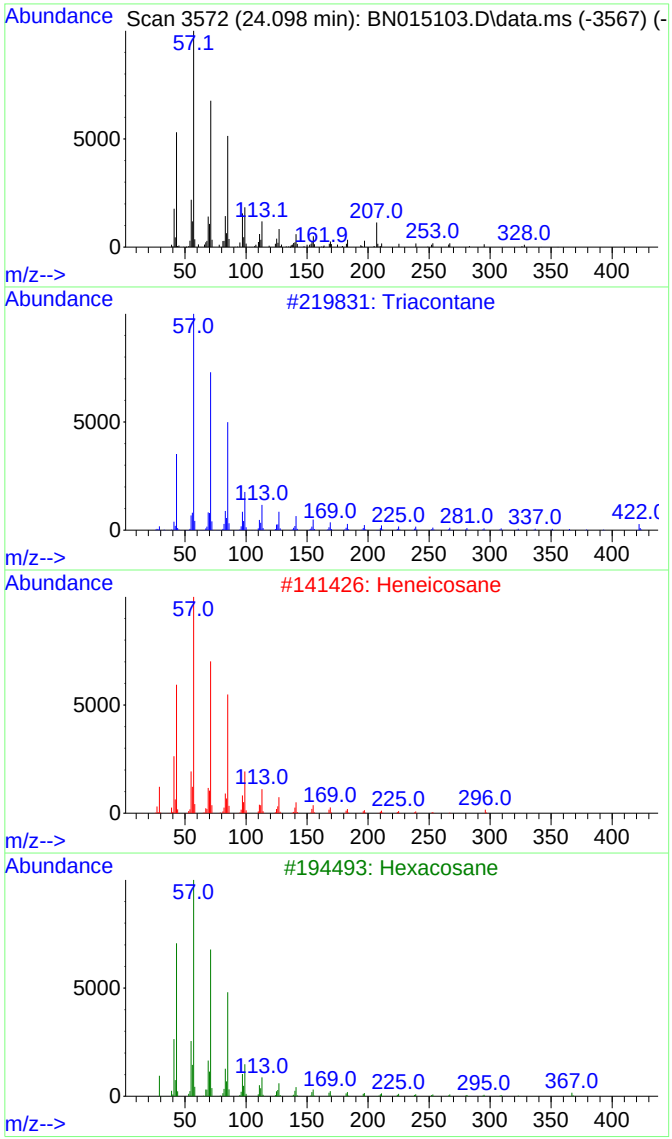
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

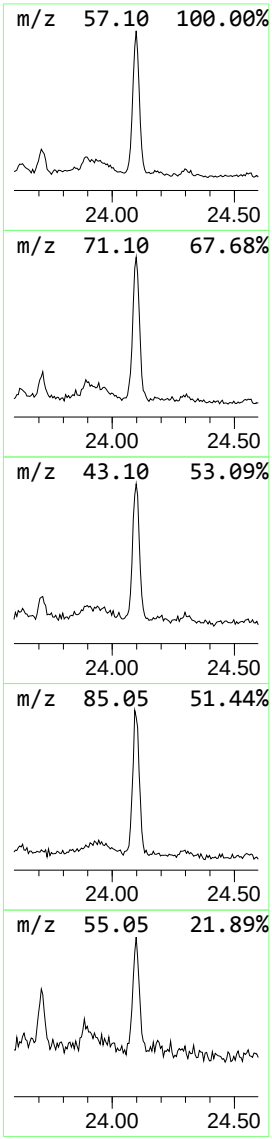
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.098	2.39 ng/ul	298915	Perylene-d12	23.539

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triacontane	422	C30H62	000638-68-6	87
2	Heneicosane	296	C21H44	000629-94-7	87
3	Hexacosane	366	C26H54	000630-01-3	87
4	2-methyloctacosane	408	C29H60	1000376-72-8	87
5	Docosane	310	C22H46	000629-97-0	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015103.D
Acq On : 26 Jun 2021 06:51
Operator : CG/JU
Sample : M2704-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDB7

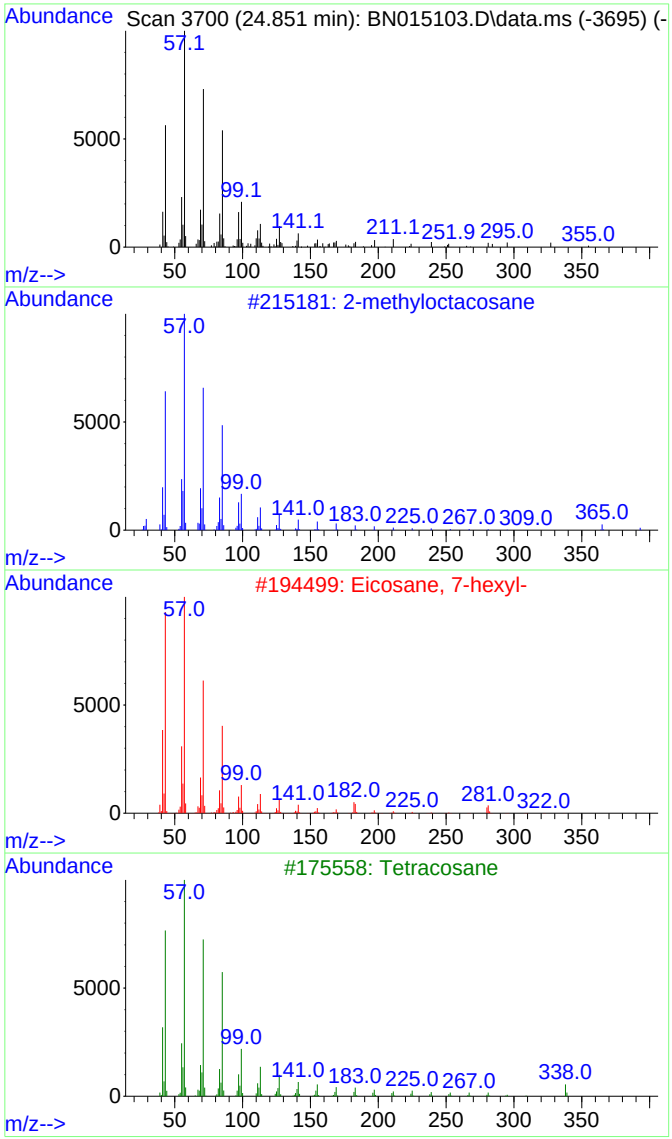
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

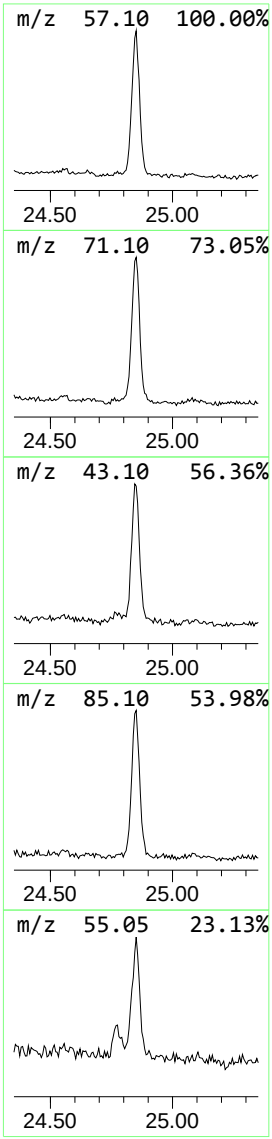
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.851	2.49 ng/ul	311938	Perylene-d12	23.539

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-methyloctacosane	408	C29H60	1000376-72-8	91
2	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
3	Tetracosane	338	C24H50	000646-31-1	90
4	Pentacosane	352	C25H52	000629-99-2	90
5	2-methylhexacosane	380	C27H56	1000376-72-7	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015103.D
Acq On : 26 Jun 2021 06:51
Operator : CG/JU
Sample : M2704-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDB7

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
p-Xylene	5.411	4.8	ng/ul	236769	1	7.728	991043	20.0
Ethanol, 2-(2-b...	10.275	3.1	ng/ul	220171	2	10.499	1418130	20.0
1,3,5-Triazine-...	15.775	2.4	ng/ul	258075	4	17.098	2122670	20.0
1,3-Benzenediol...	18.427	4.7	ng/ul	500635	4	17.098	2122670	20.0
Adipic acid, 2-...	20.516	92.1	ng/ul	10925000	5	21.292	2371890	20.0
1,3-Benzenedica...	22.157	2.1	ng/ul	244249	5	21.292	2371890	20.0
Cyclododecanol,...	22.792	2.6	ng/ul	320095	6	23.539	2503510	20.0
(DEL) Alkane: S...	23.445	2.2	ng/ul	272089	6	23.539	2503510	20.0
(DEL) Alkane: S...	24.098	2.4	ng/ul	298915	6	23.539	2503510	20.0
(DEL) Alkane: S...	24.851	2.5	ng/ul	311938	6	23.539	2503510	20.0