LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100720\
Data File : BN012096.D
Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

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\SOM-EPA-BN100220.M
Title : SVOA CALIBRATION

Signal : TIC: BN012096.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.081	13	16	24	rVB	87728	119194	0.85%	0.057%
2	3.205	32	37	47	rVB4	109841	225074	1.60%	0.108%
3	3.699	117	121	128	rVB2	47345	75614	0.54%	0.036%
4	3.822	137	142	147	rVB2	26677	44187	0.31%	0.021%
5	3.911	152	157	164	rVB	348442	489490	3.48%	0.236%
6	5.199	370	376	385	rVB	1458865	2035492	14.47%	0.980%
7	5.328	392	398	402	rBV2	36303	63845	0.45%	0.031%
8	5.669	449	456	465	rBV	9694751	13958332	99.25%	6.723%
9	6.164	534	540	546	rBV	468548	663106	4.71%	0.319%
10	6.422	578	584	588	rBV	51411	82334	0.59%	0.040%
11	6.487	588	595	597	rBV6	34418	50934	0.36%	0.025%
12	6.516	597	600	605	rVV	38040	57986	0.41%	0.028%
13	6.646	615	622	628	rVV	184587	313448	2.23%	0.151%
14	6.716	629	634	637	rVV2	101949	154655	1.10%	0.074%
15	6.752	637	640	644	rVV2	50981	78664	0.56%	0.038%
16	6.828	644	653	667	rVB2	998921	2005685	14.26%	0.966%
17	6.993	674	681	688	rBV	733331	1172642	8.34%	0.565%
18	7.046	688	690	694	rVV5	38183	51772	0.37%	0.025%
19	7.193	709	715	723	rVV	980333	1559923	11.09%	0.751%
20	7.269	723	728	739	rVV4	116679	246078	1.75%	0.119%
21	7.558	772	777	783	rVB2	35836	60211	0.43%	0.029%
22	7.652	787	793	801	rVV	1516249	2395627	17.03%	1.154%
23	7.728	801	806	813	rVV	156352	248889	1.77%	0.120%
24	8.358	906	913	920	rVV	1059531	1650119	11.73%	0.795%
25	8.469	928	932	940	rVB	115983	192812	1.37%	0.093%
26	8.746	972	979	983	rBV2	53094	85036	0.60%	0.041%
27	8.799	983	988	997	rVB	861747	1427498	10.15%	0.688%
28	9.516	1104	1110	1123	rVB	753265	1243822	8.84%	0.599%
29	9.663	1131	1135	1142	rBV2	29287	47265	0.34%	0.023%
30	9.752	1145	1150	1155	rBV4	28286	44375	0.32%	0.021%
31	10.052	1194	1201	1209	rBV	1180238	1921265	13.66%	0.925%
32	10.428	1255	1265	1269	rBV	2073363	3321410	23.62%	1.600%
33	10.475	1269	1273	1280	rVV2	883740	1424954	10.13%	0.686%
34	10.569	1282	1289	1303	rVV	585967	1112558	7.91%	0.536%
35	12.022	1533	1536	1543	rVB4	27475	47073	0.33%	0.023%
36	13.122	1719	1723	1728	rVB3	30180	46071	0.33%	0.022%

LSC Area Percent Report

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Misc :
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Instrument : BNA_N
ClientSampleId : C0AP4

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Title : SVOA CALIBRATION

37	13.193	1729	1735	1741	rBV5	36592	73941	0.53%	0.036%
38	13.704	1816	1822	1827	rVV	1956290	2752901	19.57%	1.326%
39	13.751	1827	1830	1840	rVV	919995	1271734	9.04%	0.613%
40	13.834	1840	1844	1852	rVB10	36883	69109	0.49%	0.033%
41	13.987	1861	1870	1875	rBV	2172834	3249357	23.10%	1.565%
42	14.034	1875	1878	1883	rVB	276059	382373	2.72%	0.184%
43	14.293	1915	1922	1927	rBV	2894975	4275662	30.40%	2.059%
44	14.498	1951	1957	1971	rBV	764596	1224847	8.71%	0.590%
45	14.651	1980	1983	1987	rBV3	46325	61715	0.44%	0.030%
46	14.722	1987	1995	2000	rVB3	43985	91618	0.65%	0.044%
47	14.916	2023	2028	2033	rBV9	27511	54877	0.39%	0.026%
48	15.016	2040	2045	2052	rBV2	181311	249774	1.78%	0.120%
49	15.157	2067	2069	2075	rVB2	48071	54412	0.39%	0.026%
50	15.292	2086	2092	2098	rBV	2767849	3881740	27.60%	1.870%
51	15.345	2098	2101	2106	rVB2	47701	66338	0.47%	0.032%
52	15.410	2107	2112	2120	rVV	521259	757411	5.39%	0.365%
53	15.528	2127	2132	2136	rBV4	77146	113718	0.81%	0.055%
54	15.792	2172	2177	2183	rVV7	26317	68518	0.49%	0.033%
55	15.881	2189	2192	2199	rVB3	59386	108934	0.77%	0.052%
56	16.022	2213	2216	2219	rBV	108364	144833	1.03%	0.070%
57	16.234	2248	2252	2260	rVB	341098	494091	3.51%	0.238%
58	16.351	2269	2272	2275	rBV5	30618	47553	0.34%	0.023%
59	16.398	2275	2280	2284	rVB	340381	412856	2.94%	0.199%
60	16.451	2285	2289	2297	rVB5	103355	156031	1.11%	0.075%
61	16.651	2316	2323	2328	rVB	719122	928537	6.60%	0.447%
62	16.698	2328	2331	2335	rBV3	49468	73179	0.52%	0.035%
63	16.745	2335	2339	2347	rVB	50799	99770	0.71%	0.048%
64	16.816	2347	2351	2355	rBV3	105934	157057	1.12%	0.076%
65	16.863	2355	2359	2364	rVB3	90474	142831	1.02%	0.069%
66	17.045	2383	2390	2394	rBV	3372287	4829028	34.34%	2.326%
67	17.086	2394	2397	2401	rVV	585064	739519	5.26%	0.356%
68	17.139	2401	2406	2411	rVV	2823488	4021122	28.59%	1.937%
69	17.181	2411	2413	2417	rVB2	185816	236618	1.68%	0.114%
70	17.439	2453	2457	2462	rBV2	81763	161054	1.15%	0.078%
71	17.557	2473	2477	2480	rBV3	127432	172836	1.23%	0.083%
72	17.951	2540	2544	2551	rBV2	2173504	2818783	20.04%	1.358%
73	18.110	2568	2571	2575	rBV3	132463	214235	1.52%	0.103%
74	18.245	2591	2594	2597	rBV	155962	214303	1.52%	0.103%
75	18.816	2686	2691	2699	rBV2	1805674	2351080	16.72%	1.132%
76	19.110	2737	2741	2745	rBV	853882	1138690	8.10%	0.548%

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Sample : L4184-20
Misc :
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Instrument : BNA_N
ClientSampleId : C0AP4

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
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Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Title : SVOA CALIBRATION

77	19.169	2749	2751	2756	rVB	452827	533934	3.80%	0.257%
78	19.245	2759	2764	2770	rBV	1765607	2559164	18.20%	1.233%
79	19.445	2793	2798	2802	rBV	3375395	4837425	34.40%	2.330%
80	20.010	2891	2894	2899	rVB	627300	649159	4.62%	0.313%
81	20.445	2965	2968	2974	rBV2	610399	700510	4.98%	0.337%
82	20.510	2975	2979	2983	rVB	1737100	1905786	13.55%	0.918%
83	20.727	3013	3016	3024	rVB	1154254	1465962	10.42%	0.706%
84	20.975	3055	3058	3064	rVB	5212352	6343639	45.11%	3.056%
85	21.033	3065	3068	3076	rVB	724538	934900	6.65%	0.450%
86	21.174	3088	3092	3097	rBV	1943764	2203288	15.67%	1.061%
87	21.251	3100	3105	3109	rVV	4997476	6376129	45.34%	3.071%
88	21.416	3129	3133	3138	rBV2	8535374	10099455	71.81%	4.865%
89	21.474	3139	3143	3147	rVV	3836190	4339497	30.86%	2.090%
90	21.516	3147	3150	3153	rVV	1268259	1611242	11.46%	0.776%
91	21.545	3153	3155	3164	rVB	1632875	2138789	15.21%	1.030%
92	21.845	3202	3206	3214	rBV	10081513	11255023	80.03%	5.421%
93	22.304	3279	3284	3290	rVB	9651760	12336214	87.71%	5.942%
94	22.804	3363	3369	3374	rBV	9759941	14064108	100.00%	6.774%
95	23.321	3452	3457	3459	rBV	2284111	3805957	27.06%	1.833%
96	23.357	3459	3463	3473	rVB	7720583	12428429	88.37%	5.986%
97	23.463	3476	3481	3490	rVB	2709582	4758888	33.84%	2.292%
98	23.992	3565	3571	3579	rVB	6260337	11600371	82.48%	5.588%
99	24.721	3689	3695	3703	rVB	3577683	7765387	55.21%	3.740%
100	25.580	3835	3841	3850	rVB	2418247	5817022	41.36%	2.802%

Sum of corrected areas: 207610703

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Instrument :
BNA_N
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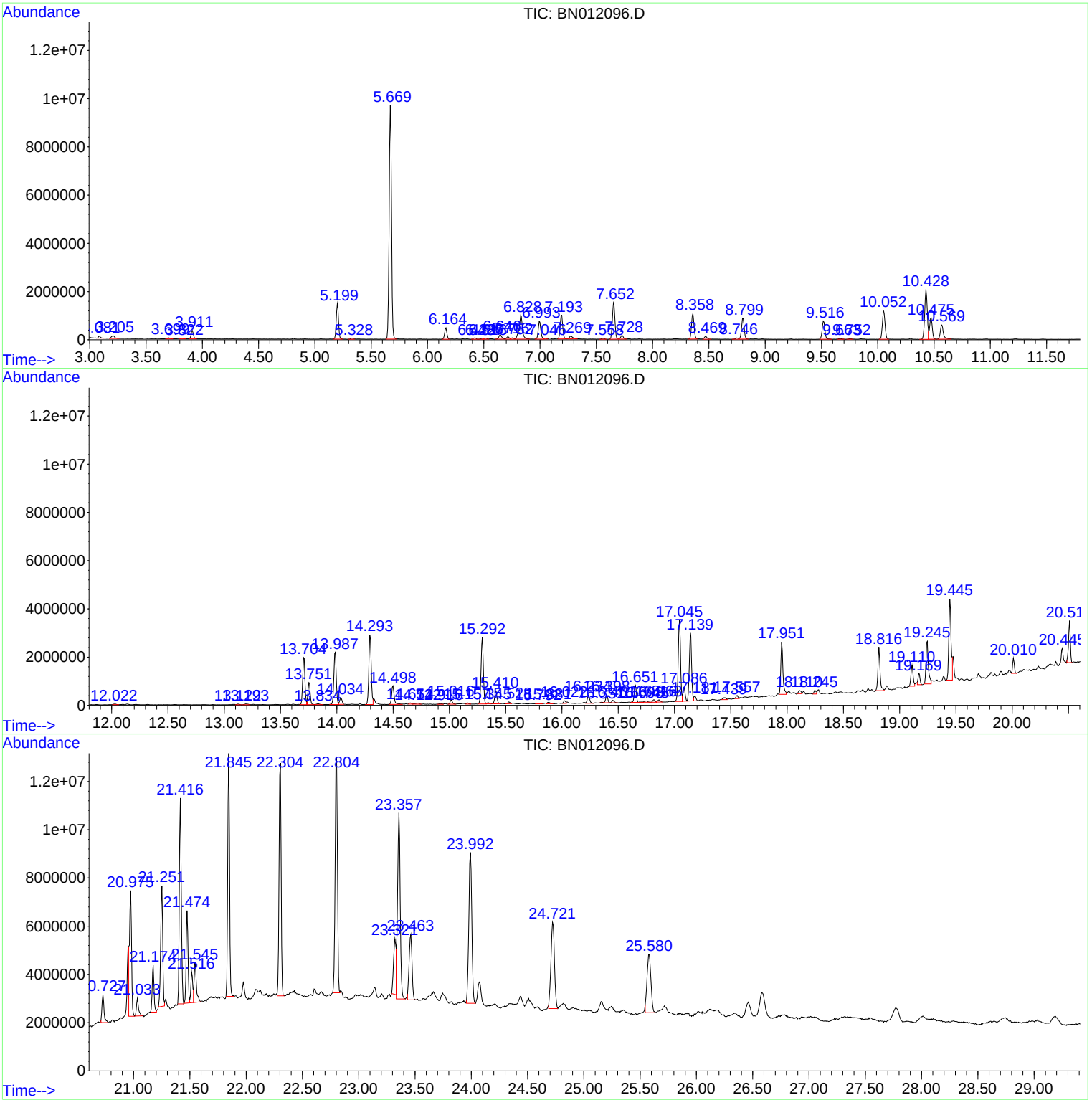
C0AP4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

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



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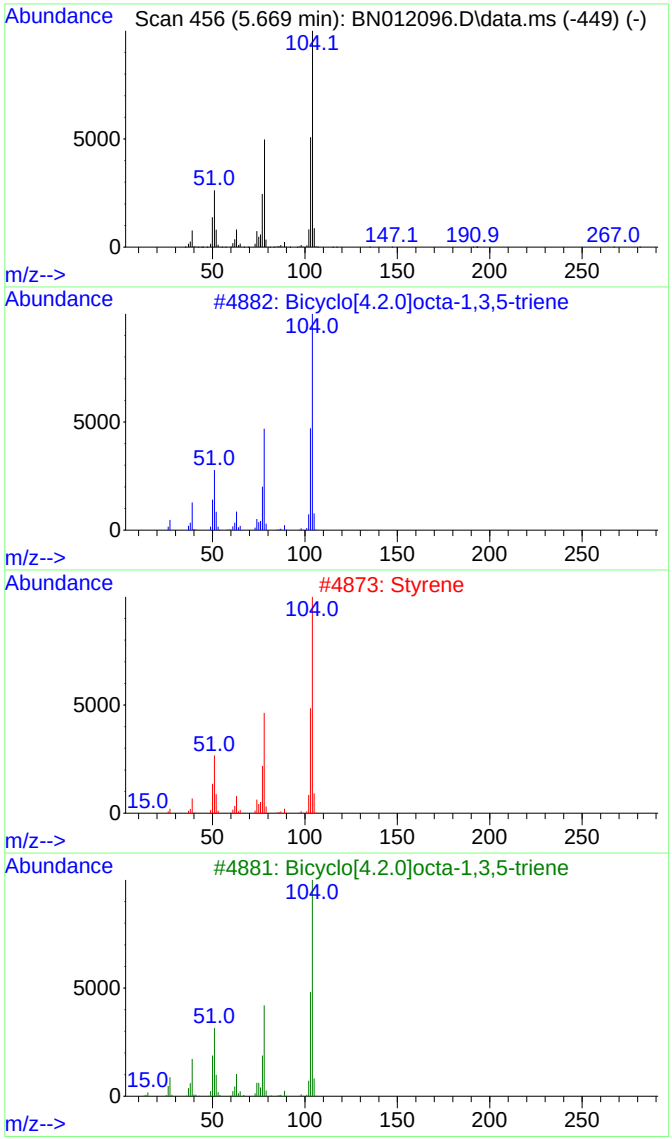
Quant Title : SVOA CALIBRATION

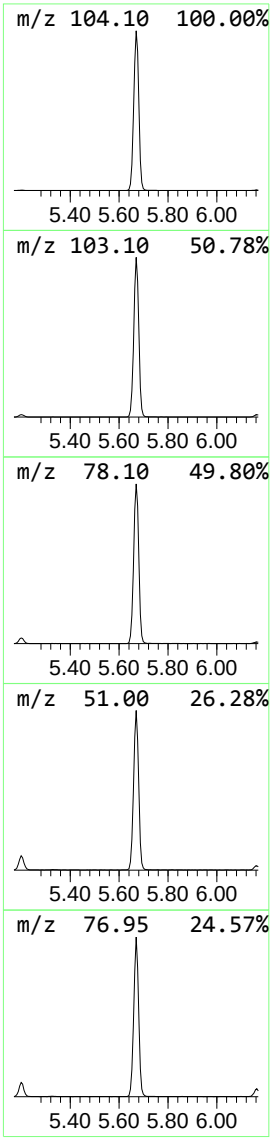
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.670	116.53 ng/ul	13958300	1,4-Dichlorobenzene-d4	7.652

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.2.0]octa-1,3,5-triene	104	C8H8	000694-87-1	96
2	Styrene	104	C8H8	000100-42-5	96
3	Bicyclo[4.2.0]octa-1,3,5-triene	104	C8H8	000694-87-1	95
4	Styrene	104	C8H8	000100-42-5	94
5	Styrene	104	C8H8	000100-42-5	93





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Data File : BN012096.D
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Instrument :
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ClientSampleId :
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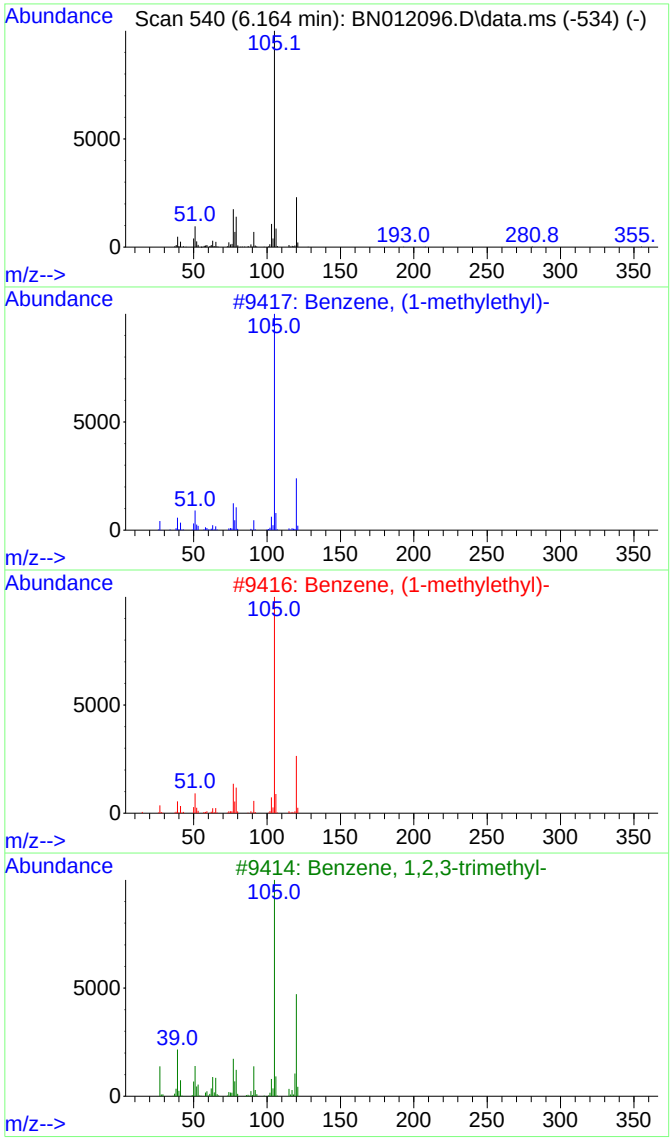
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

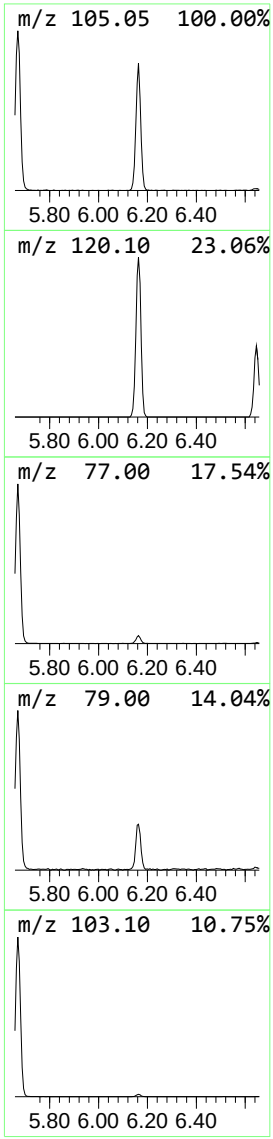
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, (1-methylethyl)- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.160	5.54 ng/ul	663106	1,4-Dichlorobenzene-d4	7.652

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	90
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	80





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Quant Title : SVOA CALIBRATION

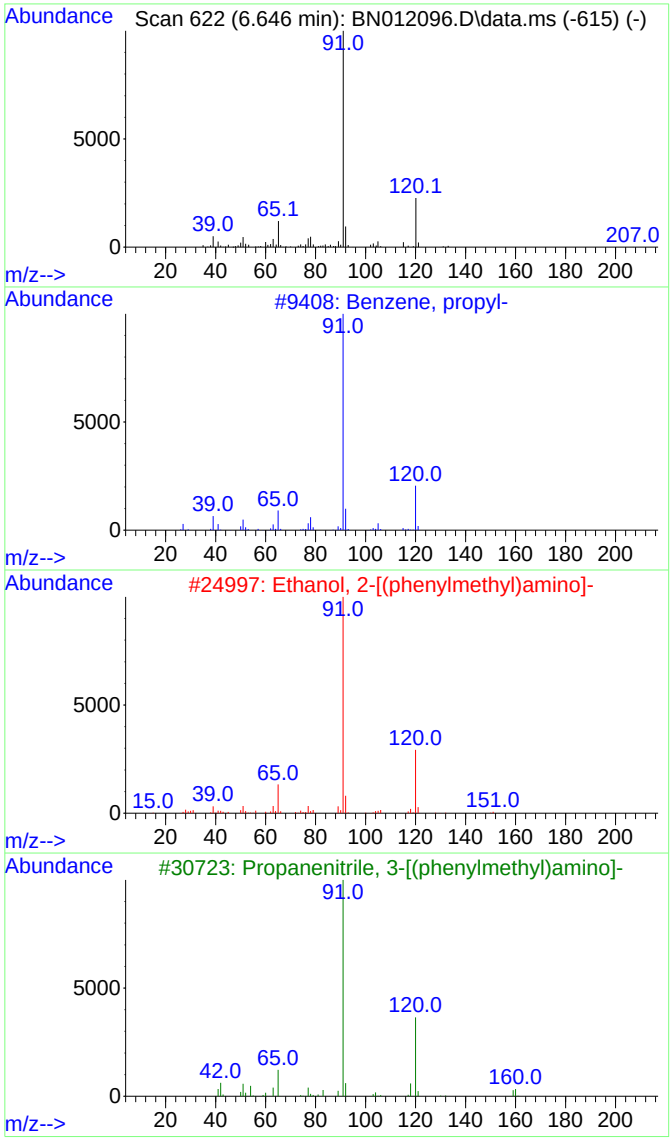
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

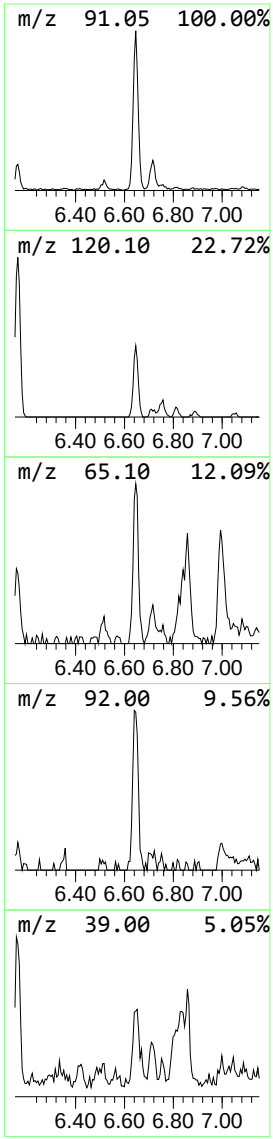
****

Peak Number 5 Benzene, propyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.650	2.62 ng/ul	313448	1,4-Dichlorobenzene-d4	7.652

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	81
2	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72
3	Propanenitrile, 3-[(phenylmethyl...	160	C10H12N2	000706-03-6	72
4	Benzene, propyl-	120	C9H12	000103-65-1	72
5	Benzeneacetaldehyde	120	C8H8O	000122-78-1	53





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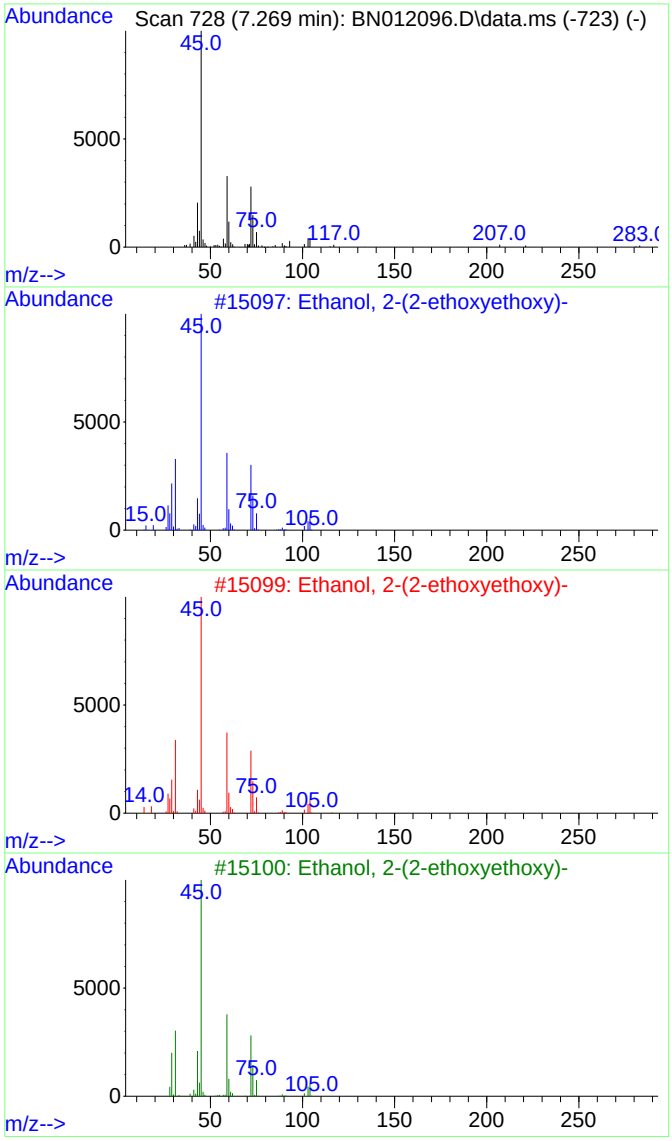
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

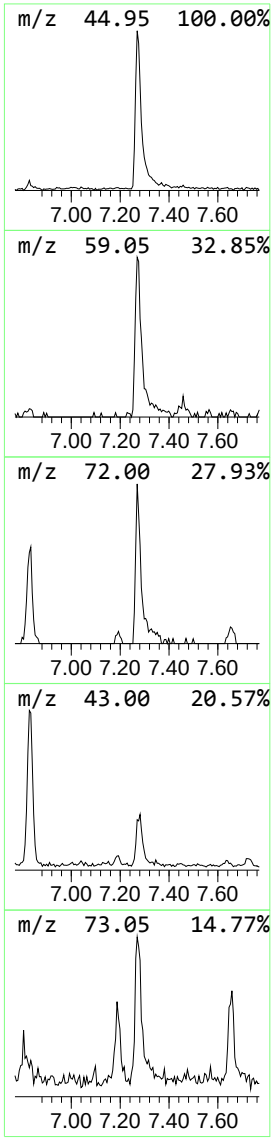
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.270	2.05 ng/ul	246078	1,4-Dichlorobenzene-d4	7.652

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
3	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	83
4	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	78
5	Diethyl carbitol	162	C8H18O3	000112-36-7	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100720\
Data File : BN012096.D
Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AP4

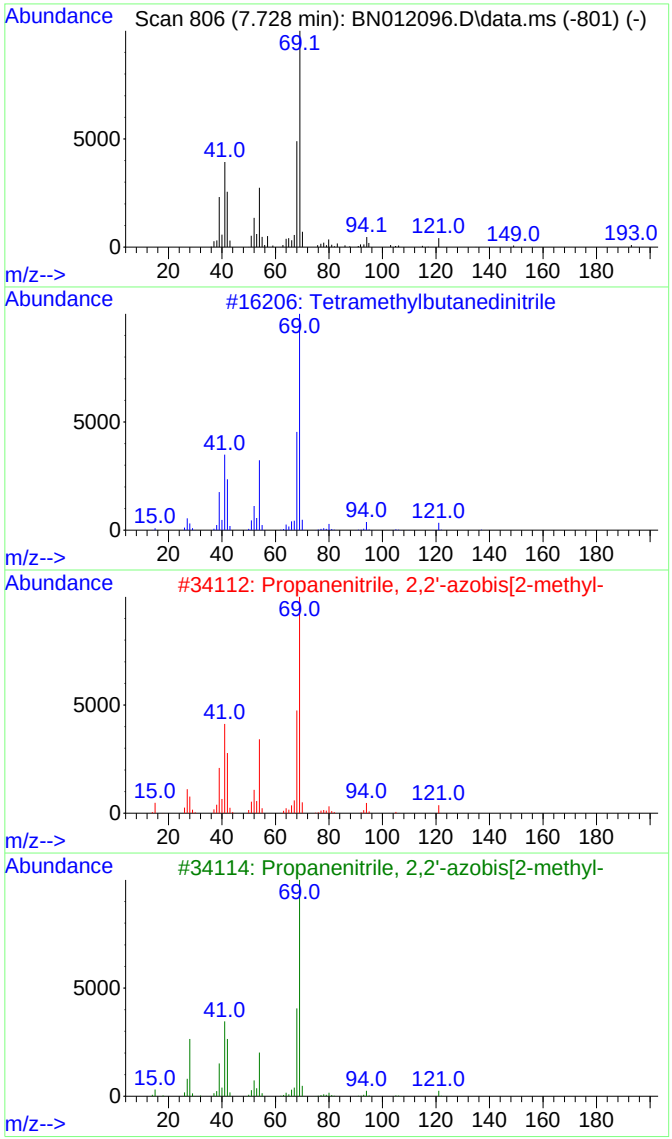
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

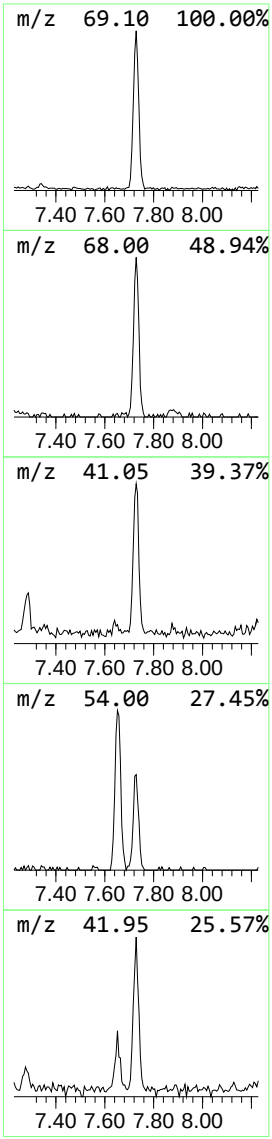
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Tetramethylbutanedinitrile Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.730	2.08 ng/ul	248889	1,4-Dichlorobenzene-d4	7.652

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetramethylbutanedinitrile	136	C8H12N2	003333-52-6	90
2	Propanenitrile, 2,2'-azobis[2-me...	164	C8H12N4	000078-67-1	90
3	Propanenitrile, 2,2'-azobis[2-me...	164	C8H12N4	000078-67-1	83
4	Tetramethylbutanedinitrile	136	C8H12N2	003333-52-6	78
5	Heptafluorobutyric acid, cyclope...	282	C9H9F7O2	1000245-81-9	50





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Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

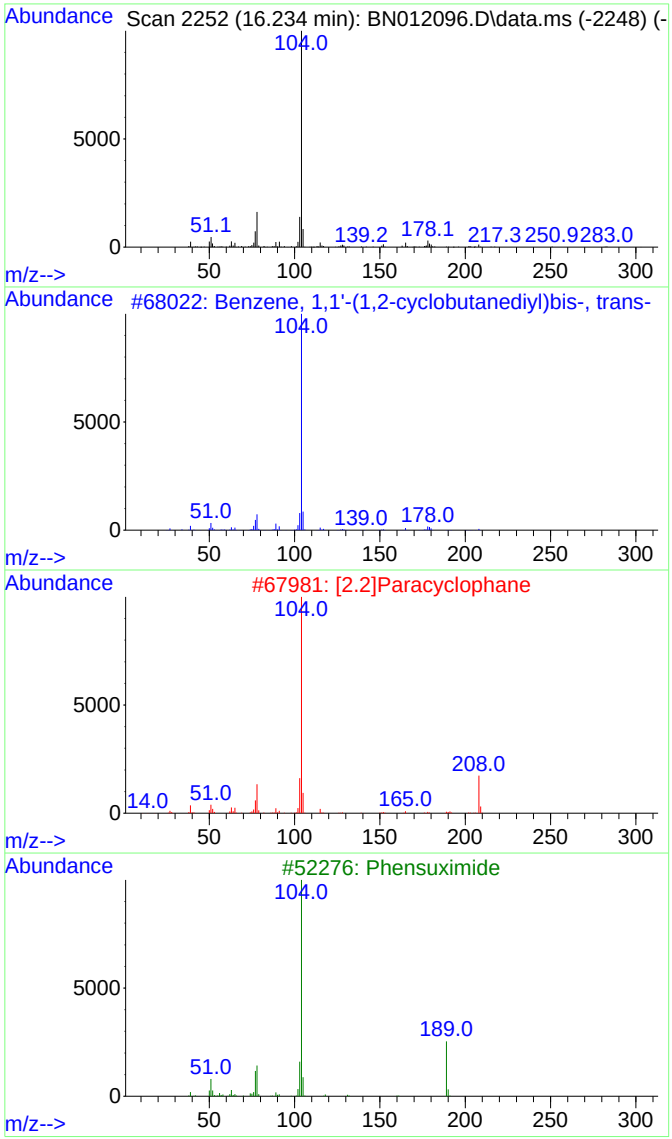
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

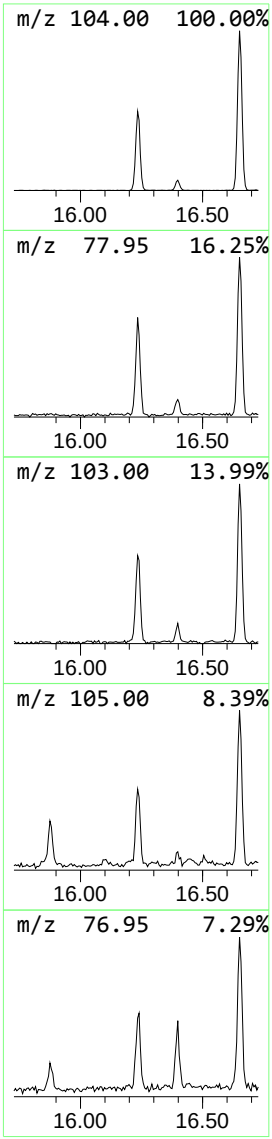
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,1'-(1,2-cyclobut... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.230	2.05 ng/ul	494091	Phenanthrene-d10	17.045

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	87
2	[2.2]Paracyclophane	208	C16H16	001633-22-3	87
3	Phensuximide	189	C11H11NO2	000086-34-0	83
4	Cyclobutane, 1,2-diphenyl-	208	C16H16	003018-21-1	78
5	[2.2]Paracyclophane	208	C16H16	001633-22-3	72





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Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

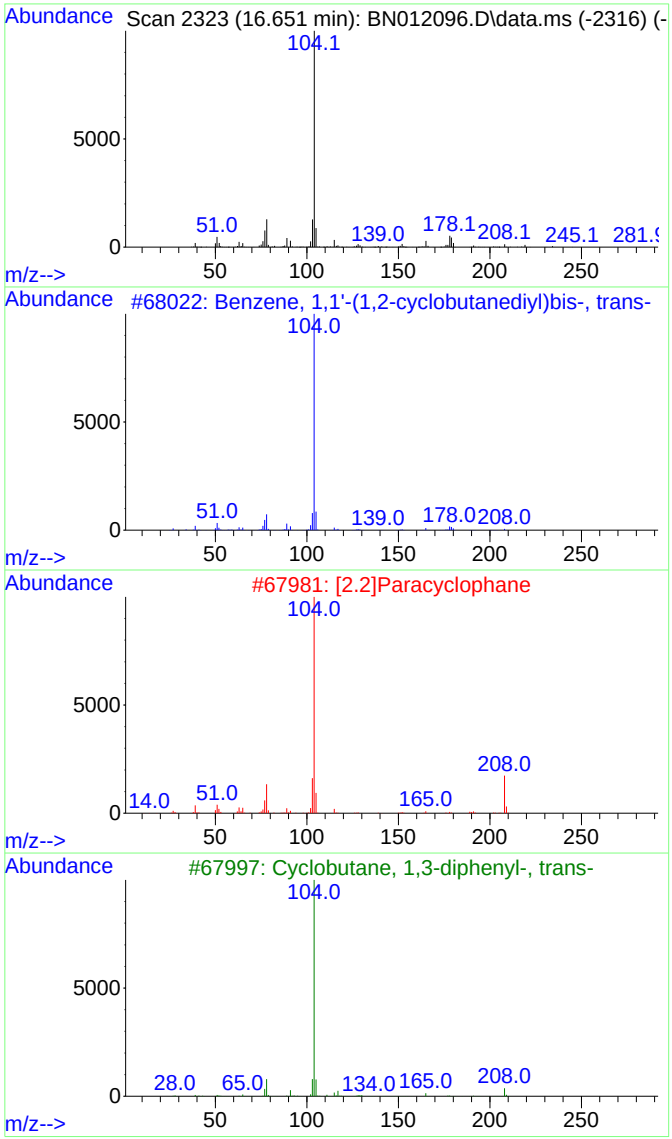
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

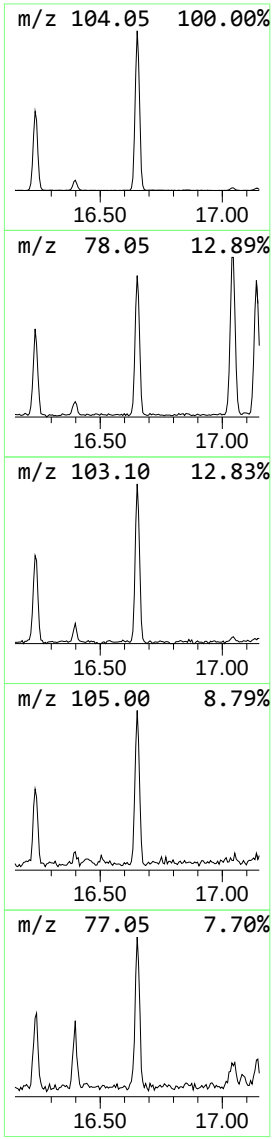
****

Peak Number 9 [2.2]Paracyclophane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.650	3.85 ng/ul	928537	Phenanthrene-d10	17.045

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	90
2	[2.2]Paracyclophane	208	C16H16	001633-22-3	74
3	Cyclobutane, 1,3-diphenyl-, trans-	208	C16H16	025558-23-0	64
4	[2.2]Paracyclophane	208	C16H16	001633-22-3	64
5	Styrene	104	C8H8	000100-42-5	58





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Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Title : SVOA CALIBRATION

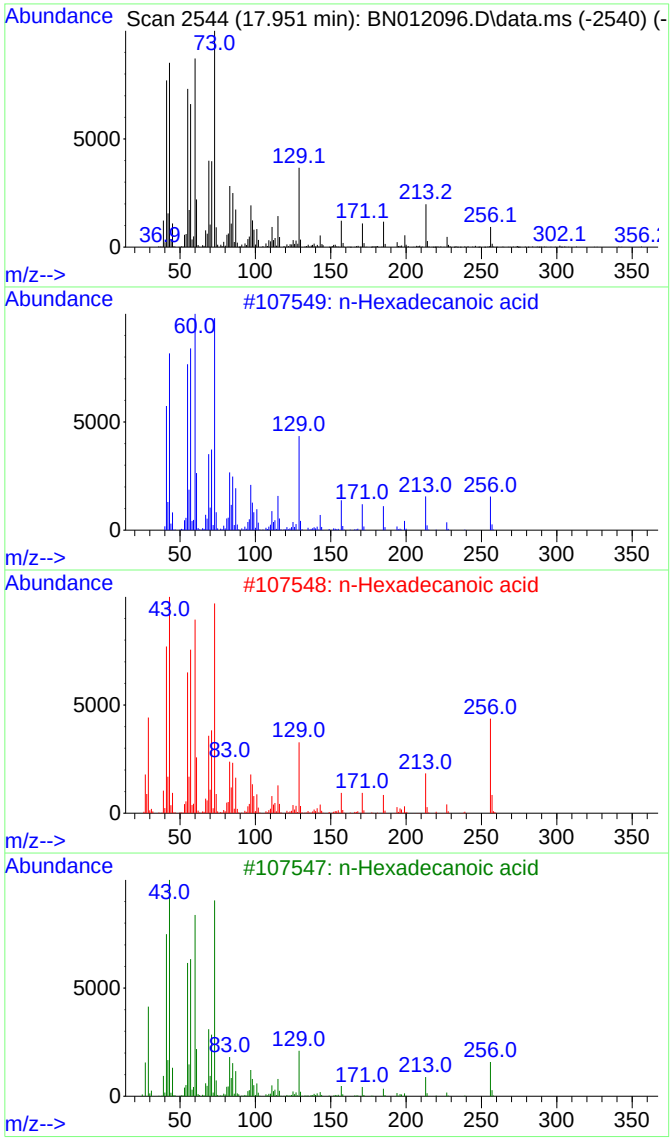
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

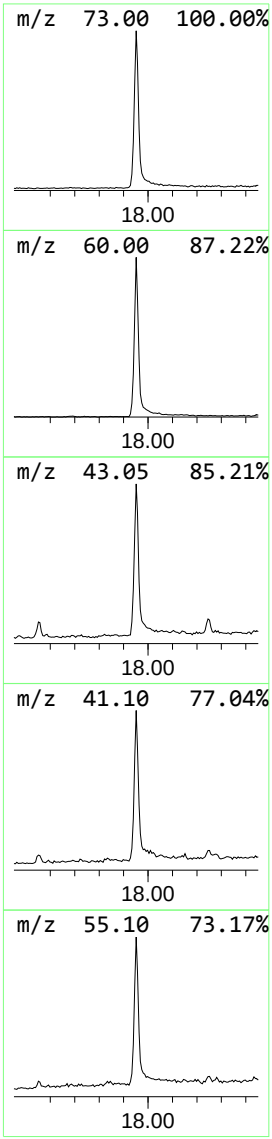
****

Peak Number 10 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.950	11.67 ng/ul	2818780	Phenanthrene-d10	17.045

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
4	Tridecanoic acid	214	C13H26O2	000638-53-9	92
5	Octadecanoic acid	284	C18H36O2	000057-11-4	87





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Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AP4

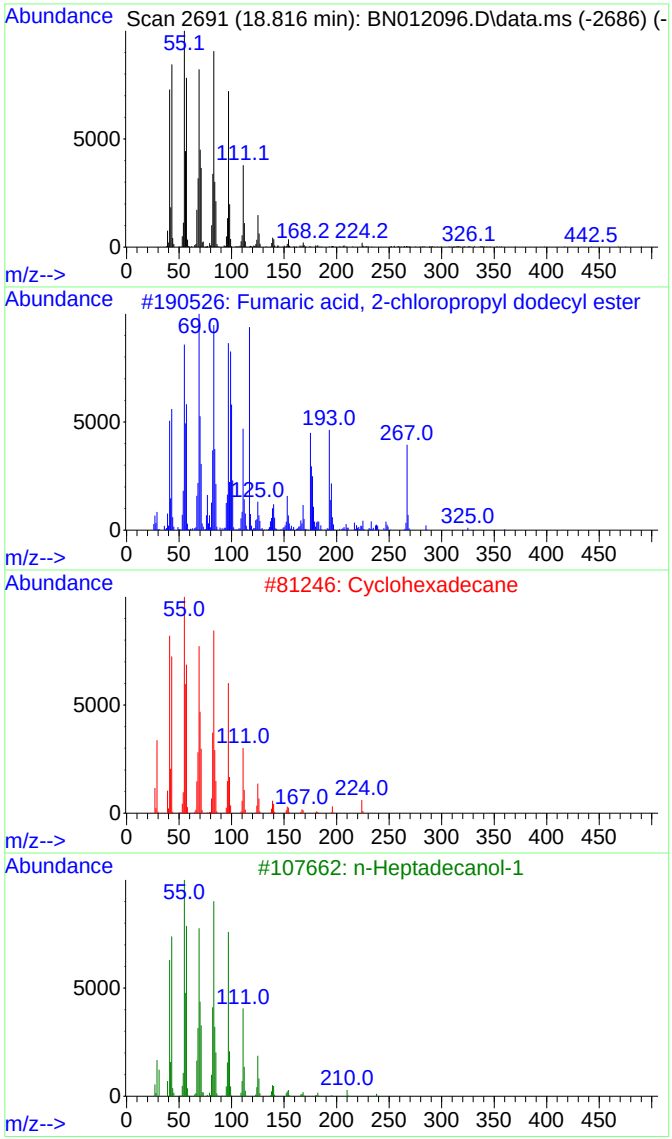
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

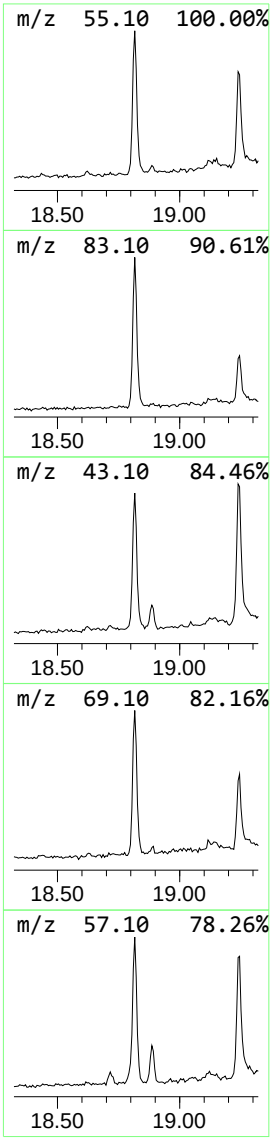
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Fumaric acid, 2-chloropropy... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.820	9.74 ng/ul	2351080	Phenanthrene-d10	17.045

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fumaric acid, 2-chloropropyl dod...	360	C19H33ClO4	1000348-57-0	98
2	Cyclohexadecane	224	C16H32	000295-65-8	95
3	n-Heptadecanol-1	256	C17H36O	001454-85-9	95
4	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
5	Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	94





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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
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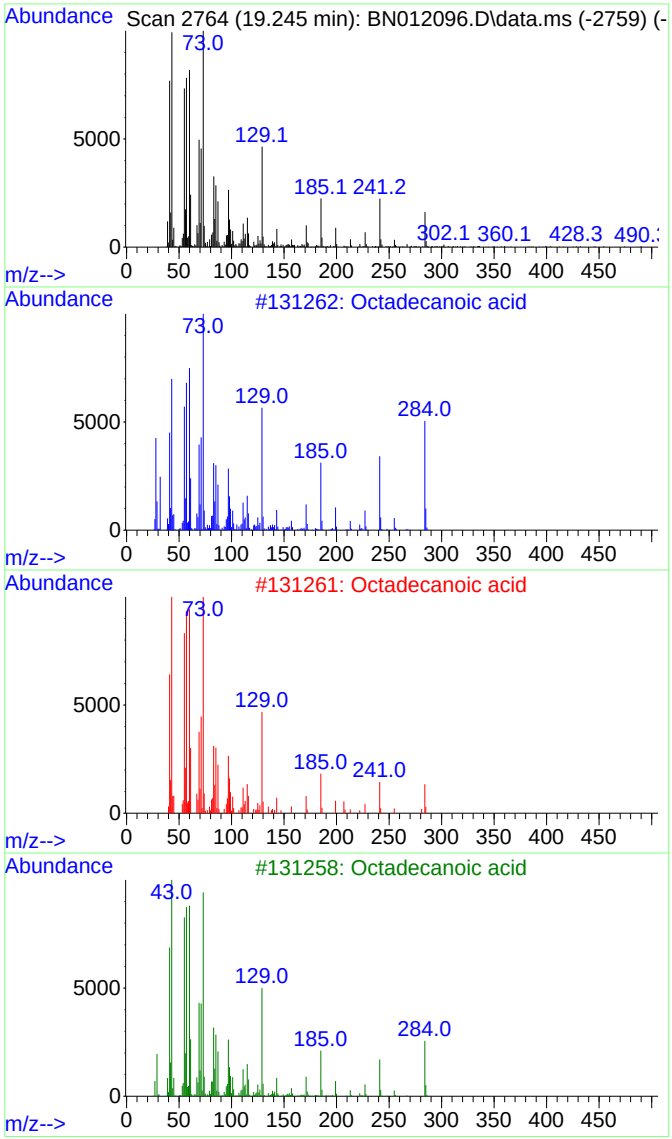
Quant Title : SVOA CALIBRATION

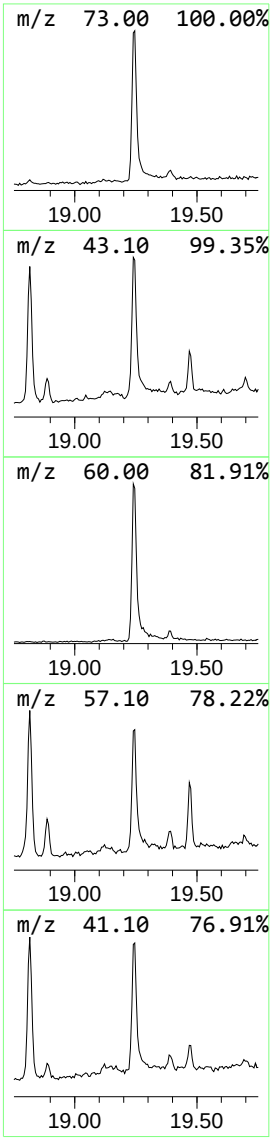
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Octadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.250	8.03 ng/ul	2559160	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	99
4	Octadecanoic acid	284	C18H36O2	000057-11-4	97
5	Octadecanoic acid	284	C18H36O2	000057-11-4	93





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Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

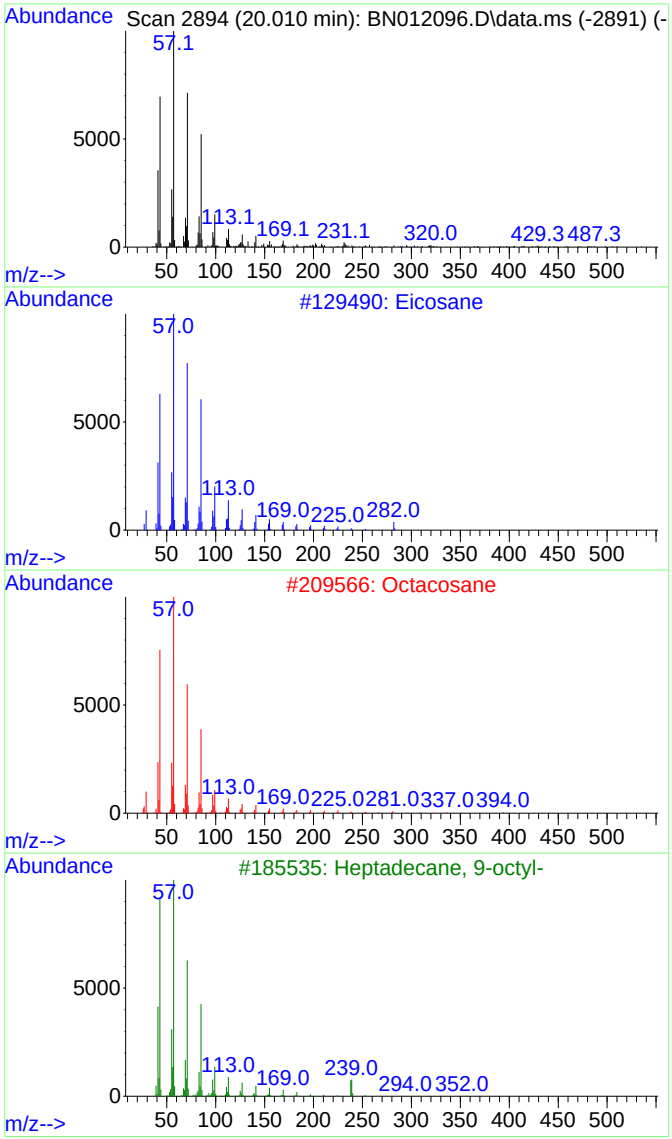
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

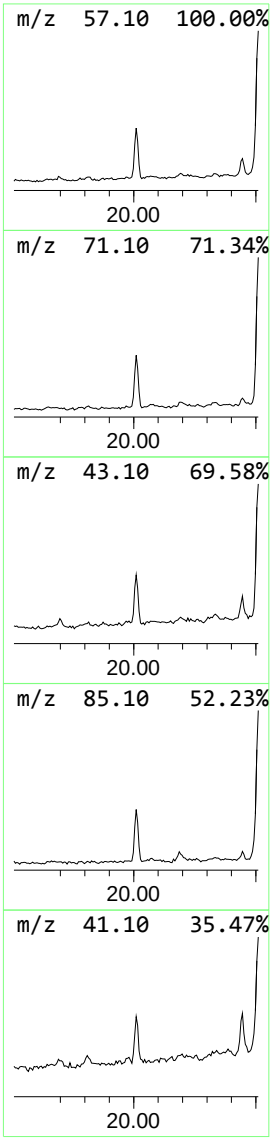
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.010	2.04 ng/ul	649159	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	95
2	Octacosane	394	C28H58	000630-02-4	93
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4	Heneicosane	296	C21H44	000629-94-7	87
5	Boric acid, ethyl-, didecyl ester	354	C22H47BO2	1000152-34-4	83





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Data File : BN012096.D
Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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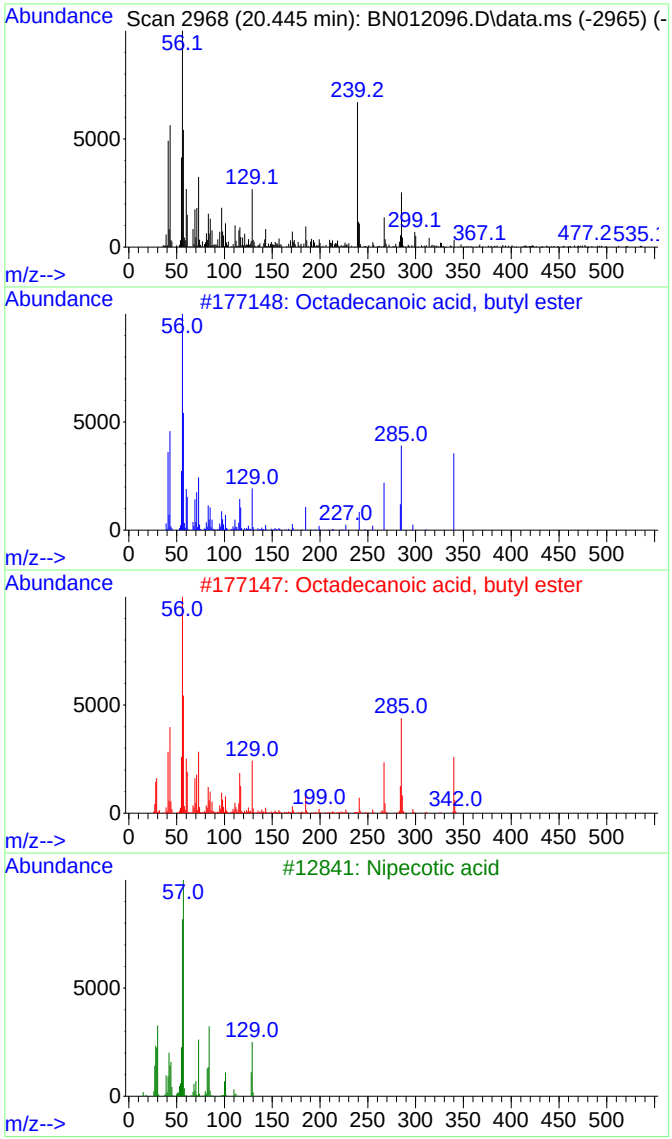
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

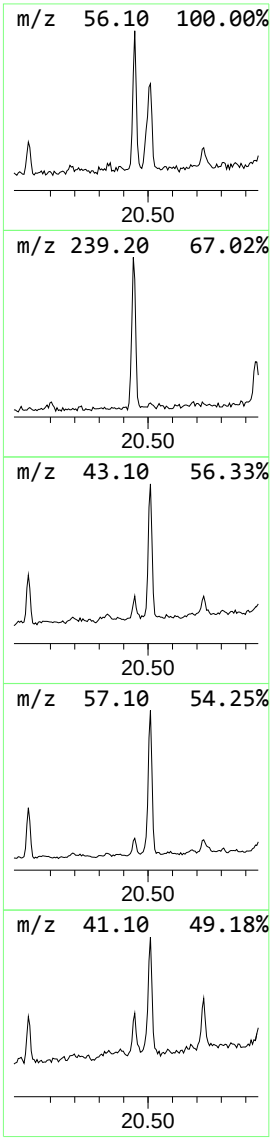
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Octadecanoic acid, butyl ester Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.450	2.20 ng/ul	700510	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	86
2	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	49
3	Nipecotic acid	129	C6H11NO2	000498-95-3	43
4	Cyclopropane, 1-(1-methylethyl)-...	210	C15H30	041977-39-3	25
5	Docosanoic acid	340	C22H44O2	000112-85-6	25





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Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

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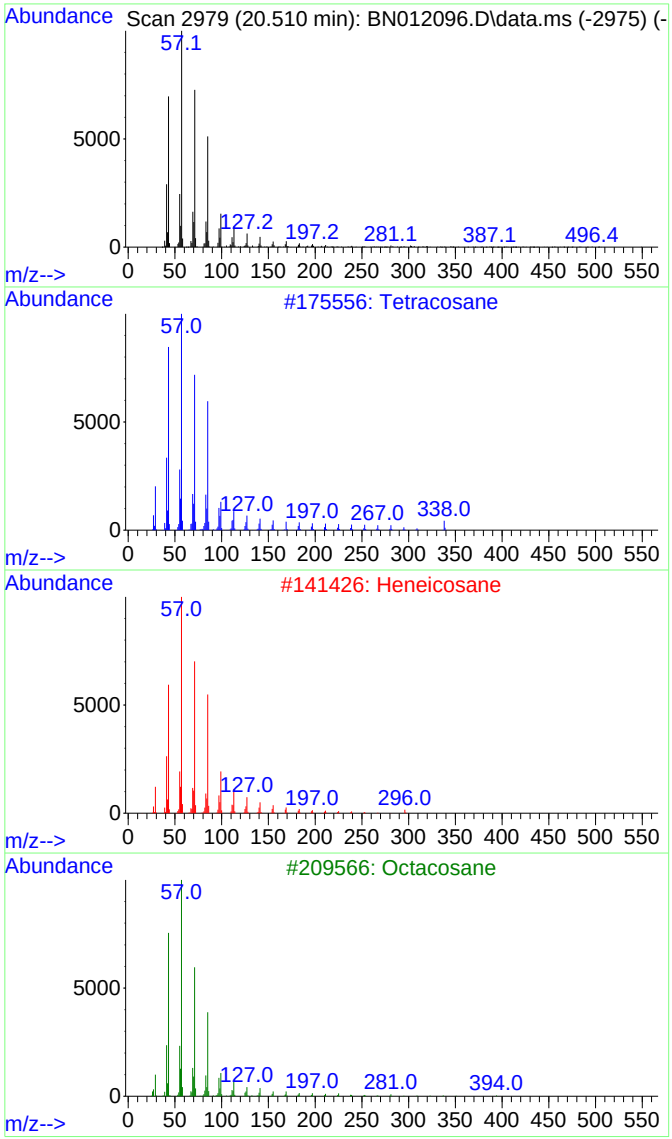
Quant Title : SVOA CALIBRATION

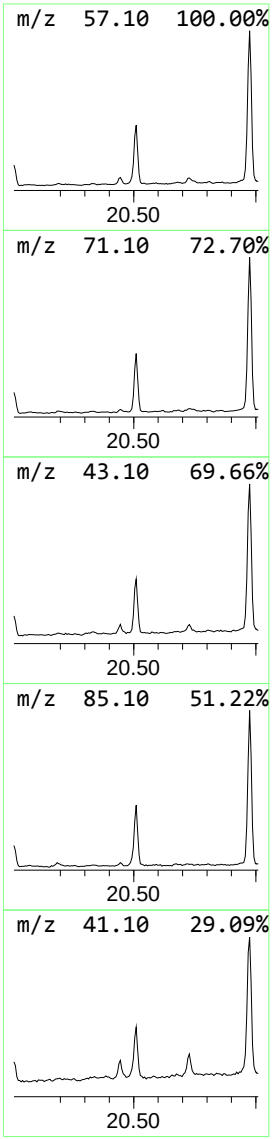
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.510	5.98 ng/ul	1905790	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	93
2	Heneicosane	296	C21H44	000629-94-7	91
3	Octacosane	394	C28H58	000630-02-4	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	90





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Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

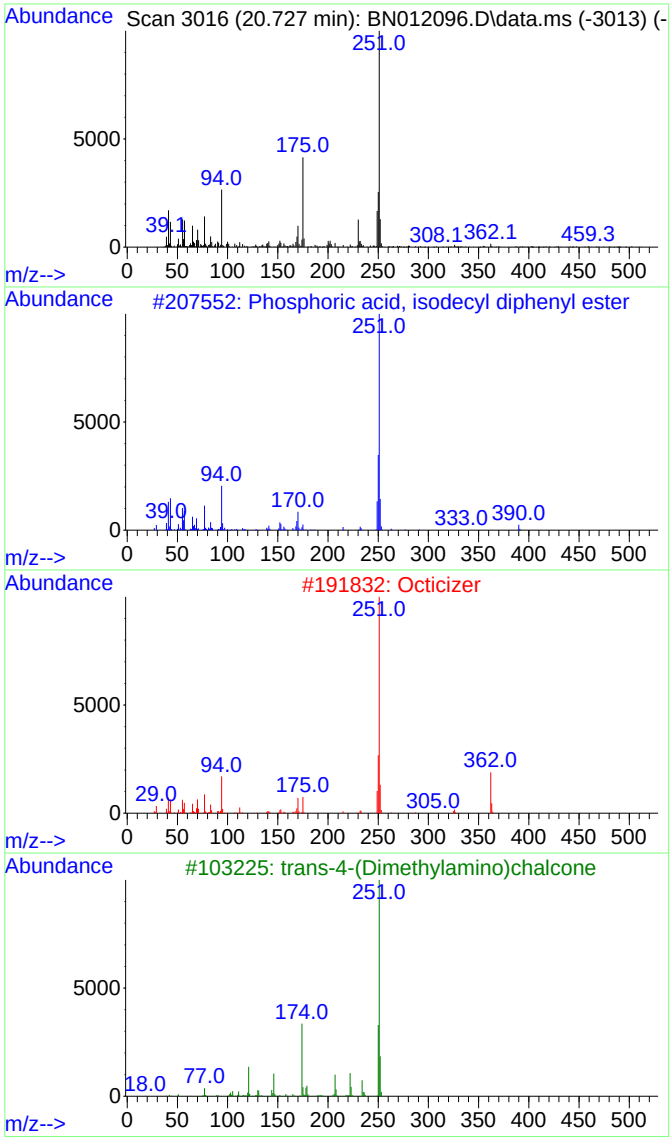
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

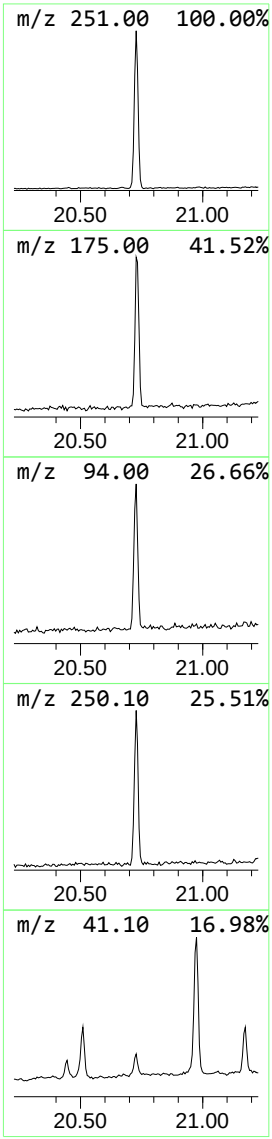
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Phosphoric acid, isodecyl d... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.730	4.60 ng/ul	1465960	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phosphoric acid, isodecyl diphen...	390	C22H31O4P	029761-21-5	58
2	Octicizer	362	C20H27O4P	001241-94-7	53
3	trans-4-(Dimethylamino)chalcone	251	C17H17NO	022965-98-6	43
4	2H-1,3,4-Benzotriazepin-2-one, 1...	251	C15H13N3O	002855-57-4	38
5	7H-Pyrazolo[4,3-E][1,2,4]triazol...	251	C12H9N7	1000310-01-6	32





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Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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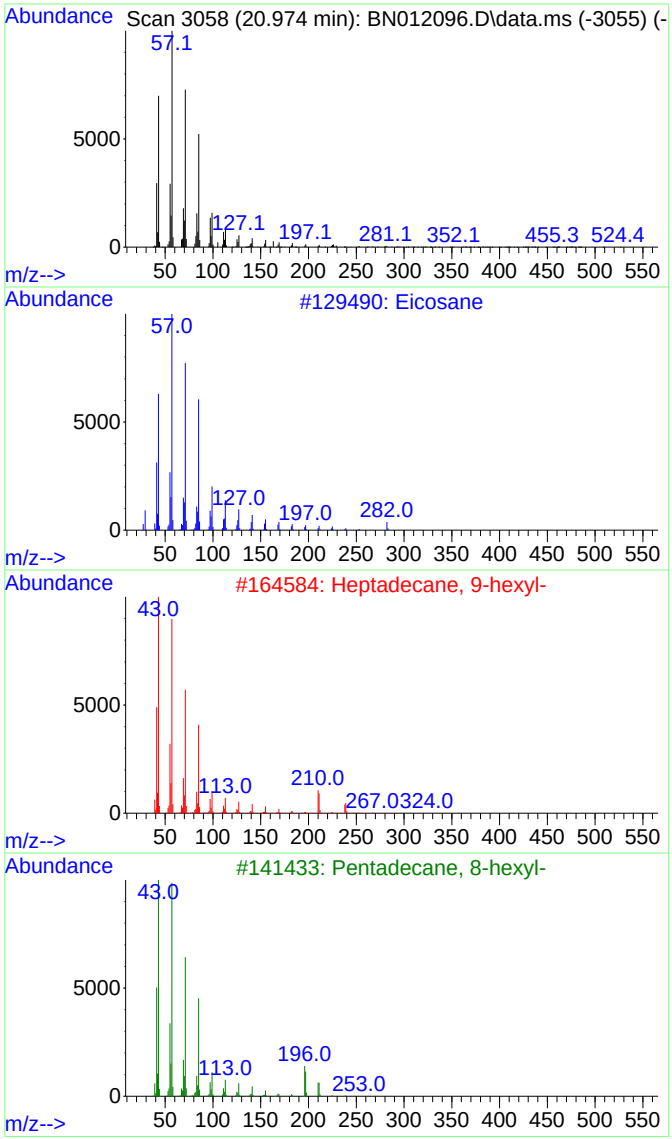
Quant Title : SVOA CALIBRATION

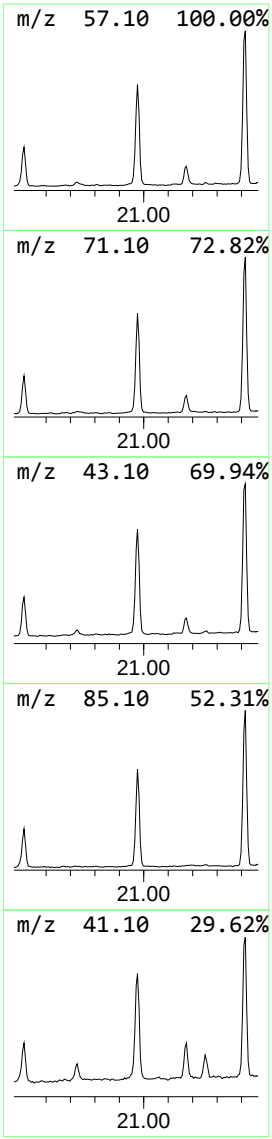
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.970	19.90 ng/ul	6343640	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	95
2	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
3	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5	Hexacosane	366	C26H54	000630-01-3	91





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Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

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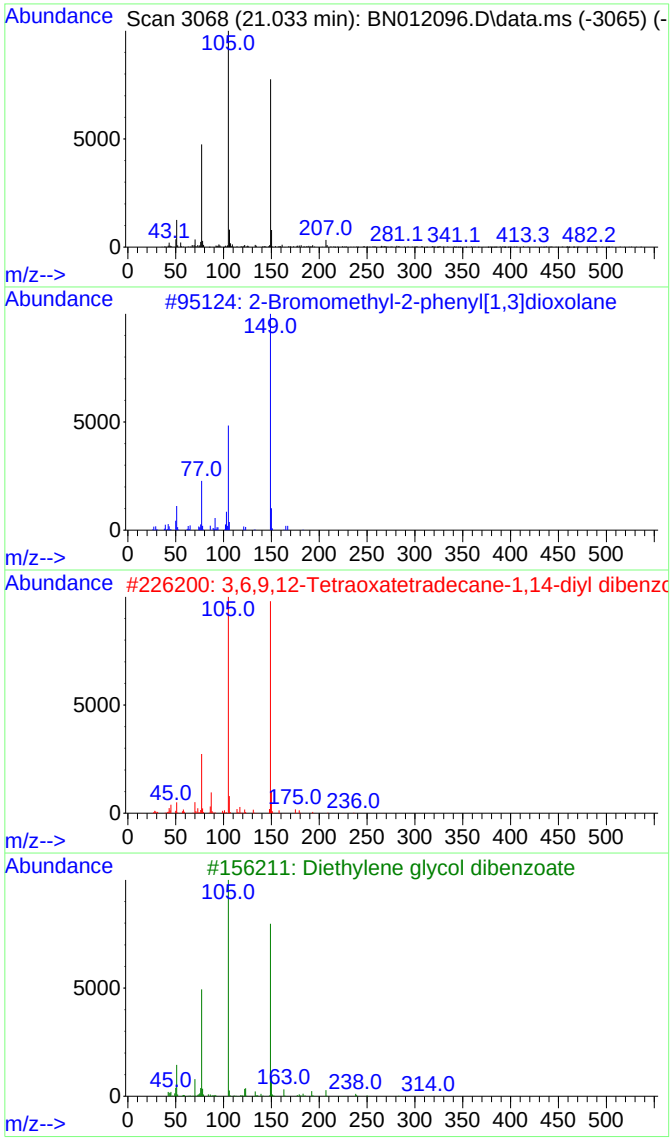
Quant Title : SVOA CALIBRATION

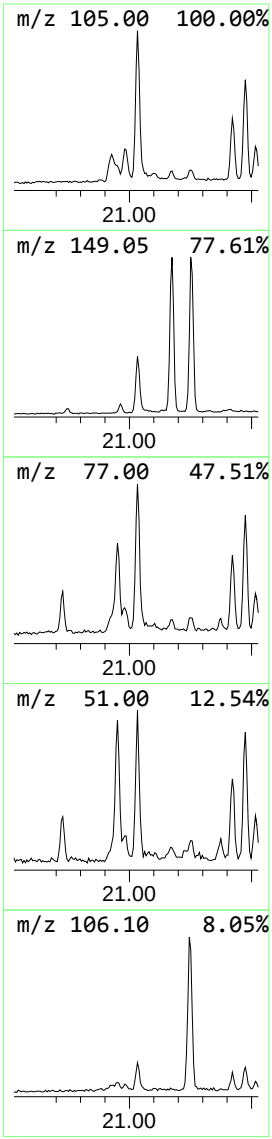
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 2-Bromomethyl-2-phenyl[1,3]... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.030	2.93 ng/ul	934900	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromomethyl-2-phenyl[1,3]dioxo...	242	C10H11BrO2	003418-21-1	64
2	3,6,9,12-Tetraoxatetradecane-1,1...	446	C24H30O8	1000366-97-0	64
3	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	64
4	Benzoic acid, 2-(4-nitrophenoxy)...	287	C15H13NO5	1000368-92-7	56
5	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	53





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Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

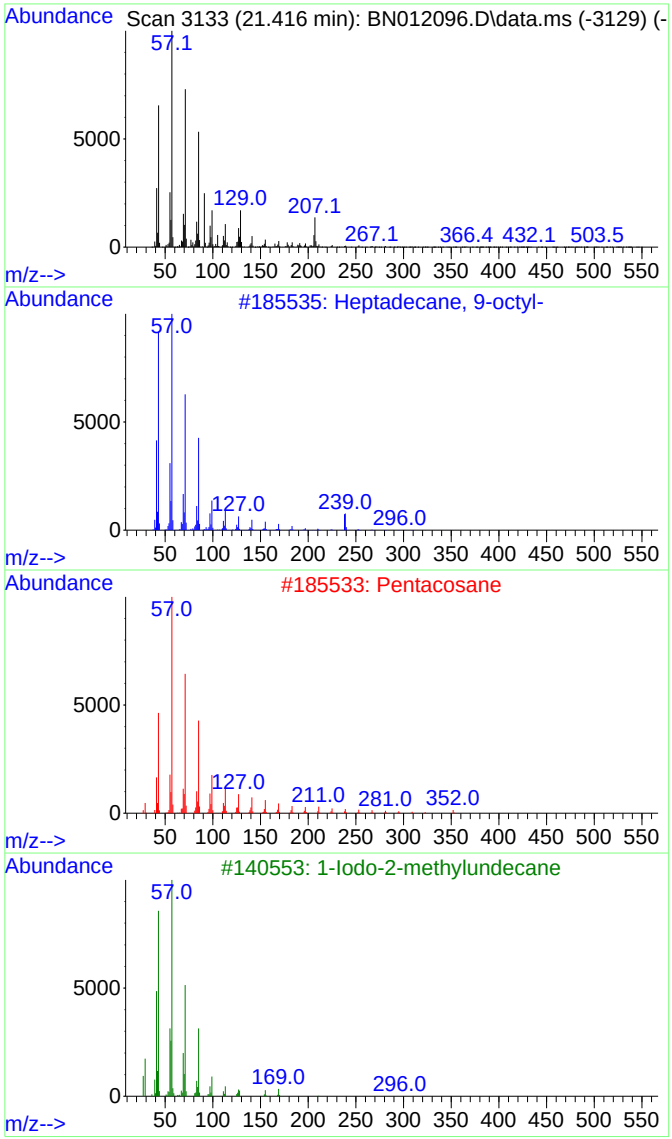
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

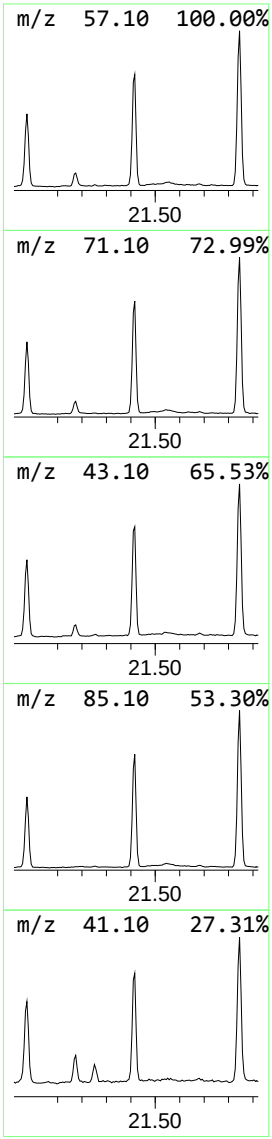
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.420	31.68 ng/ul	10099500	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2	Pentacosane	352	C25H52	000629-99-2	94
3	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	90
4	Heneicosane	296	C21H44	000629-94-7	81
5	Octacosane	394	C28H58	000630-02-4	81





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Data File : BN012096.D
Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

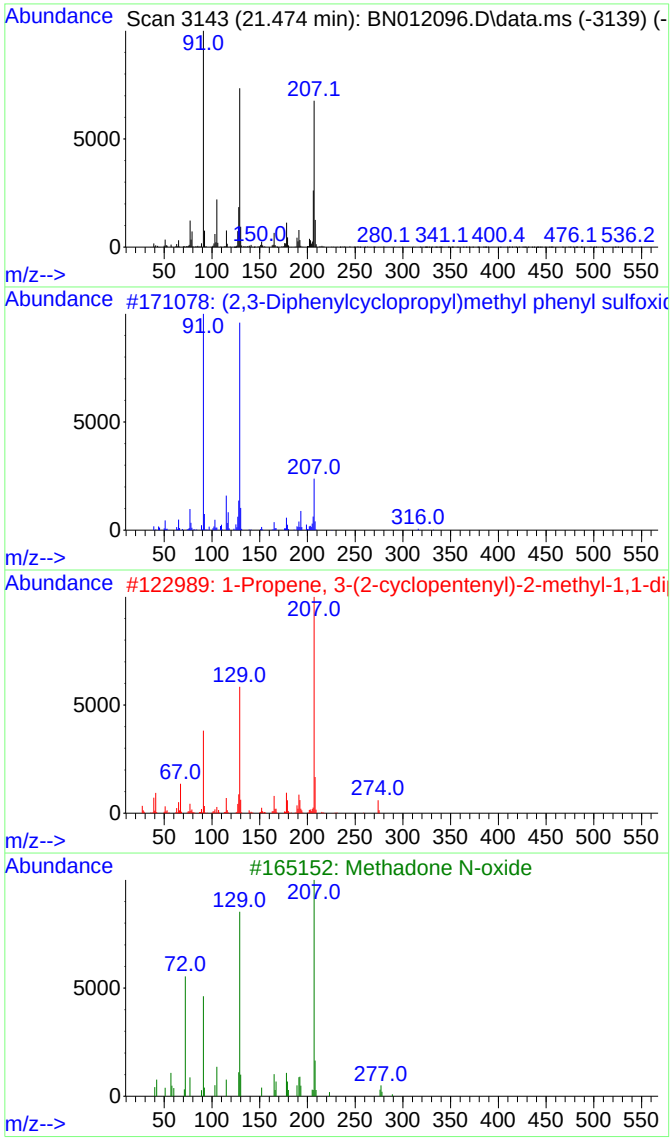
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

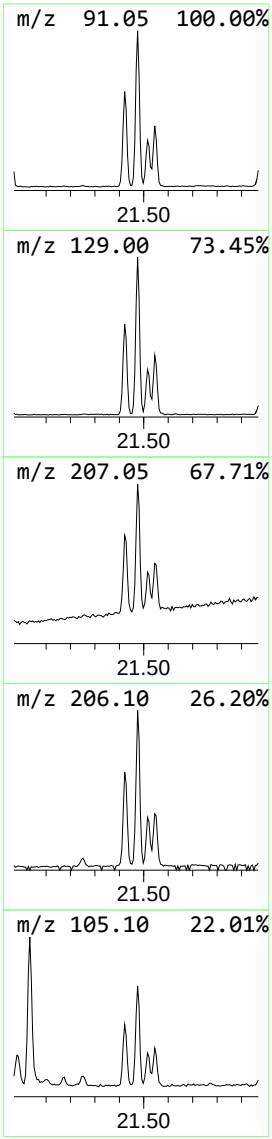
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 (2,3-Diphenylcyclopropyl)me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.470	13.61 ng/ul	4339500	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	72
2	1-Propene, 3-(2-cyclopentenyl)-2...	274	C21H22	1000154-23-3	59
3	Methadone N-oxide	325	C21H27NO2	1000120-80-7	58
4	Cyclopentene-1-carboxylic acid, ...	332	C23H24O2	1000159-40-6	40
5	Quinoline	129	C9H7N	000091-22-5	38





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Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

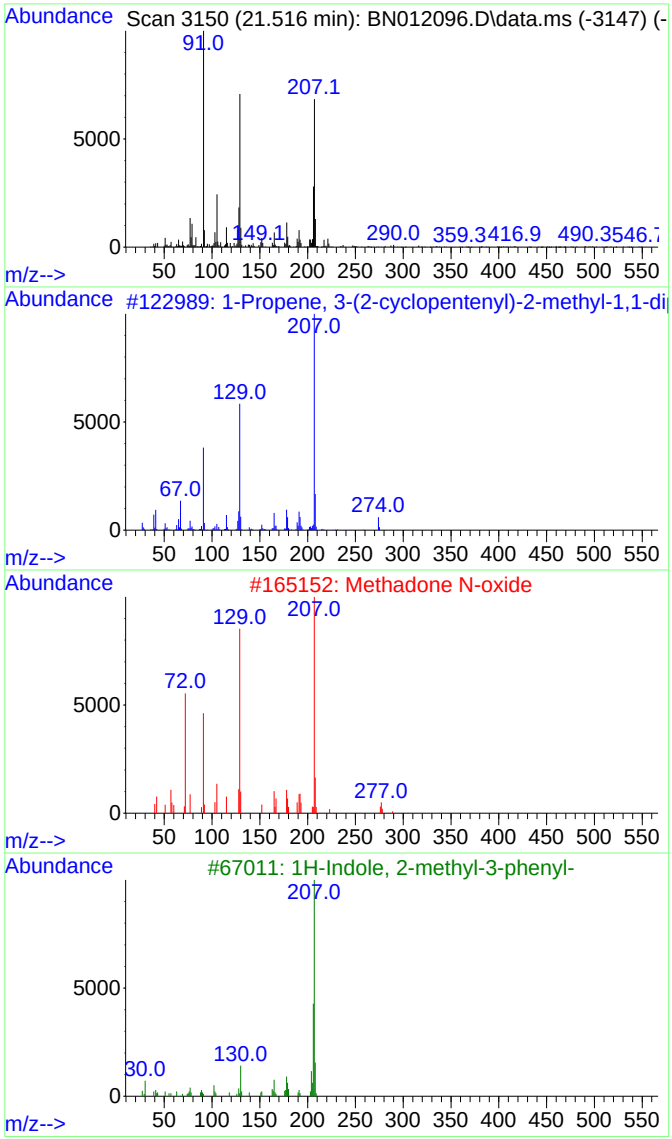
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

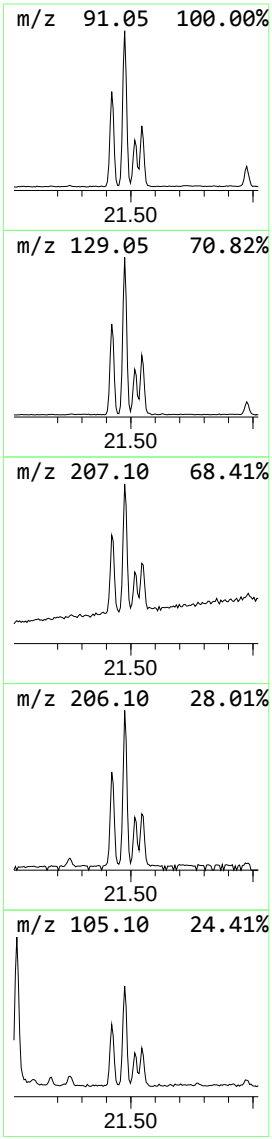
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 1-Propene, 3-(2-cyclopenten... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.520	5.05 ng/ul	1611240	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propene, 3-(2-cyclopentenyl)-2...	274	C21H22	1000154-23-3	53
2	Methadone N-oxide	325	C21H27NO2	1000120-80-7	50
3	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	41
4	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	38
5	5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	35





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Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AP4

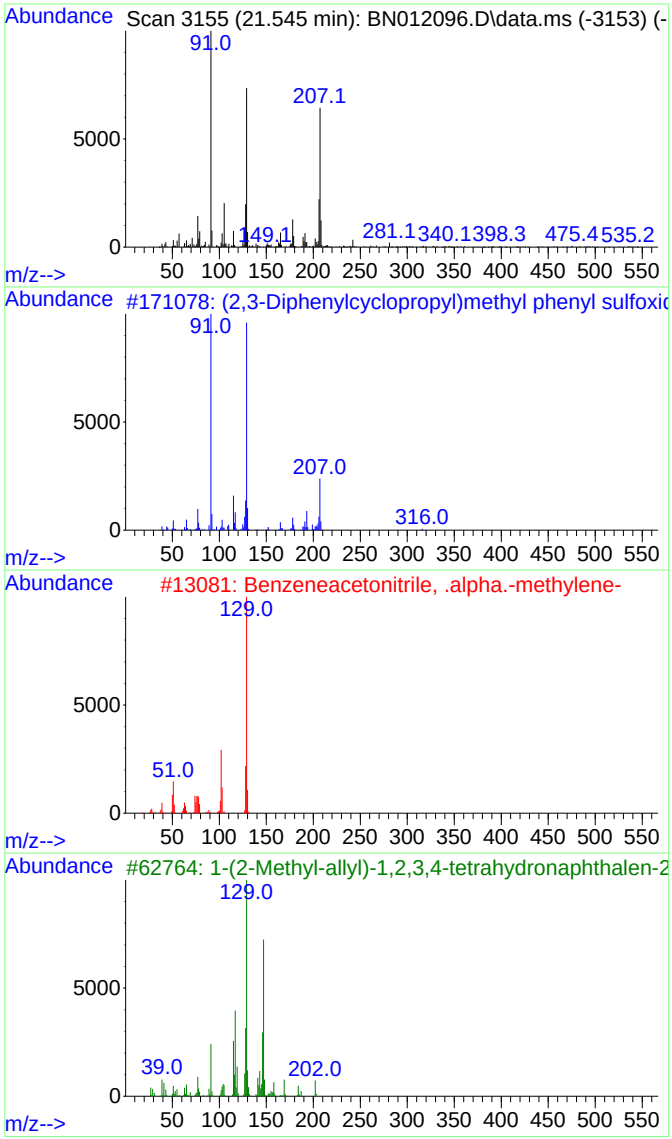
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

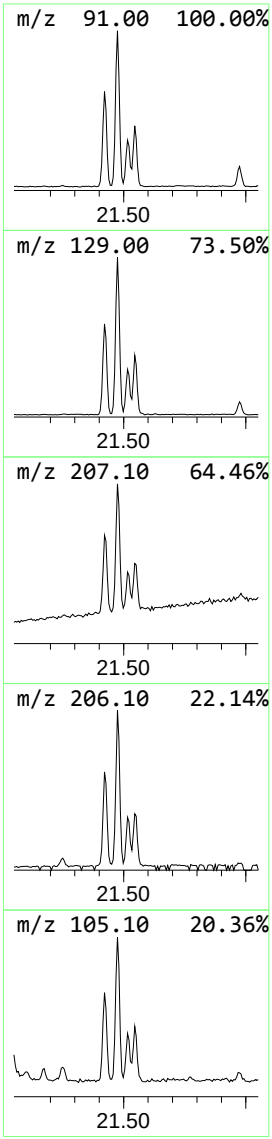
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 unknown-01 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.550	6.71 ng/ul	2138790	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	49
2	Benzeneacetonitrile, .alpha.-met...	129	C9H7N	000495-10-3	35
3	1-(2-Methyl-allyl)-1,2,3,4-tetra...	202	C14H18O	1000192-52-9	35
4	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	35
5	1-Propene, 3-(2-cyclopentenyl)-2...	274	C21H22	1000154-23-3	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100720\
Data File : BN012096.D
Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

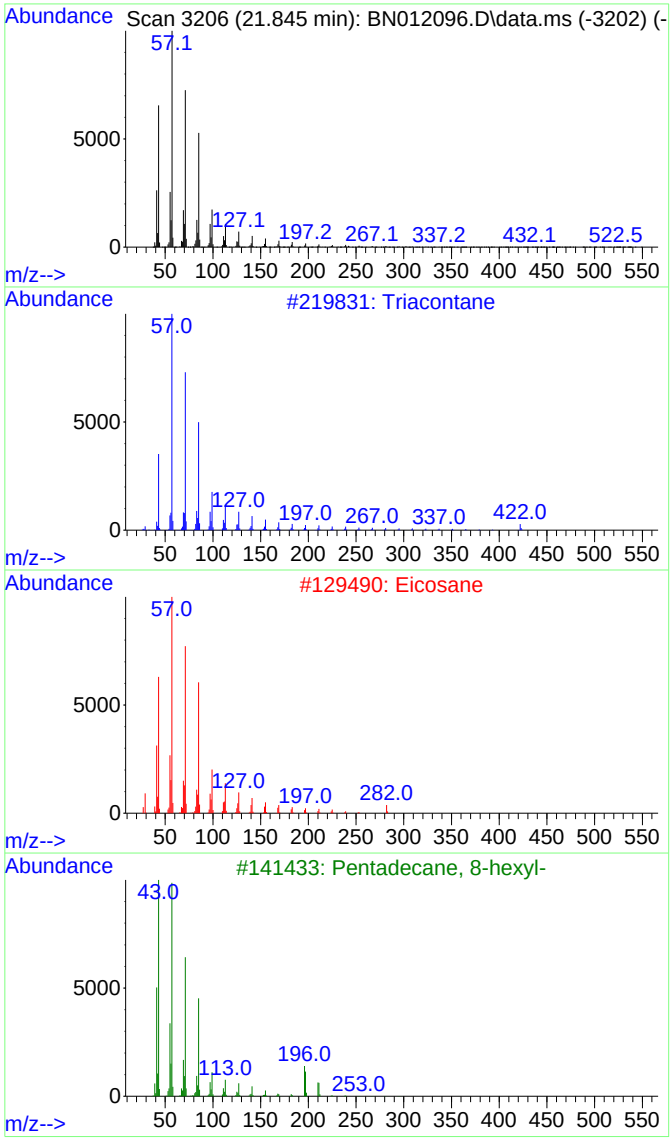
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

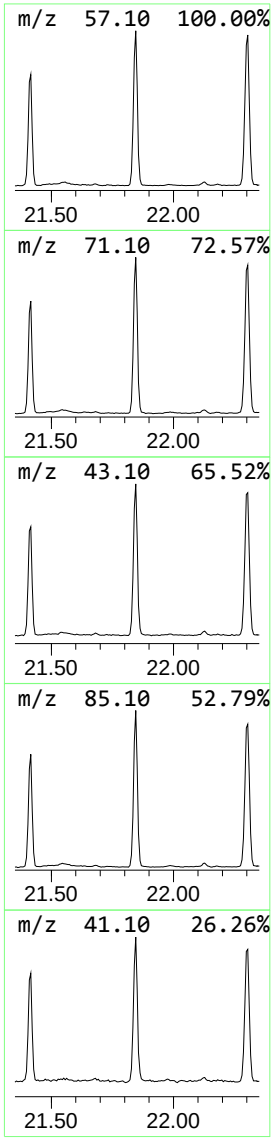
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.850	35.30 ng/ul	11255000	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triacontane	422	C30H62	000638-68-6	98
2	Eicosane	282	C20H42	000112-95-8	95
3	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
4	Octacosane	394	C28H58	000630-02-4	94
5	Tetracosane	338	C24H50	000646-31-1	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100720\
Data File : BN012096.D
Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

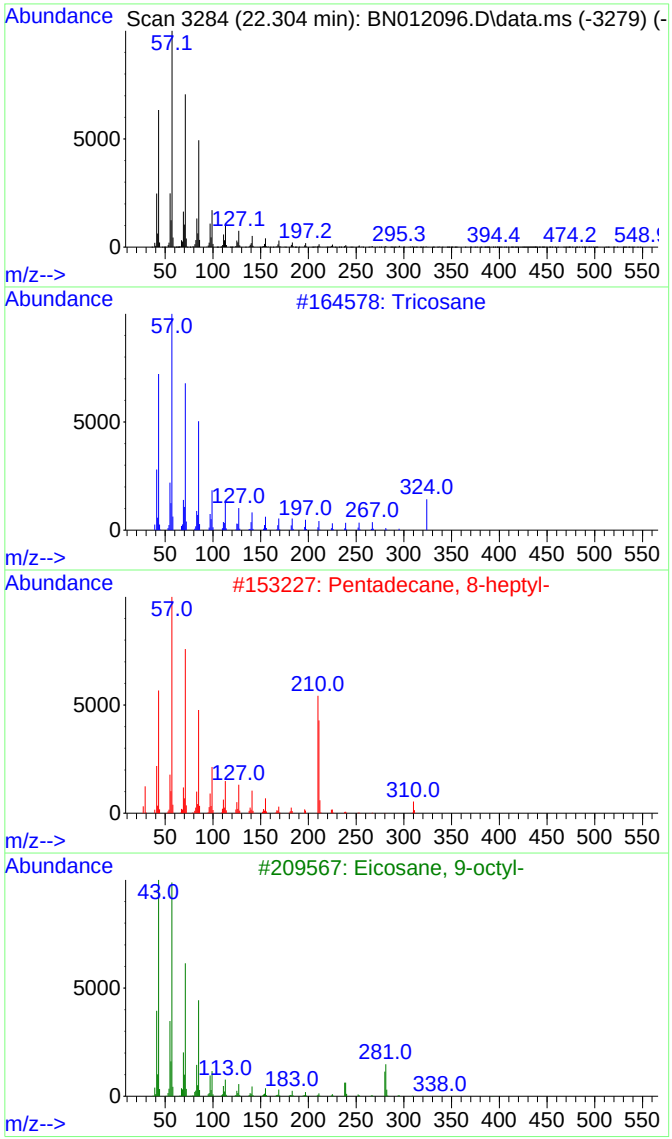
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

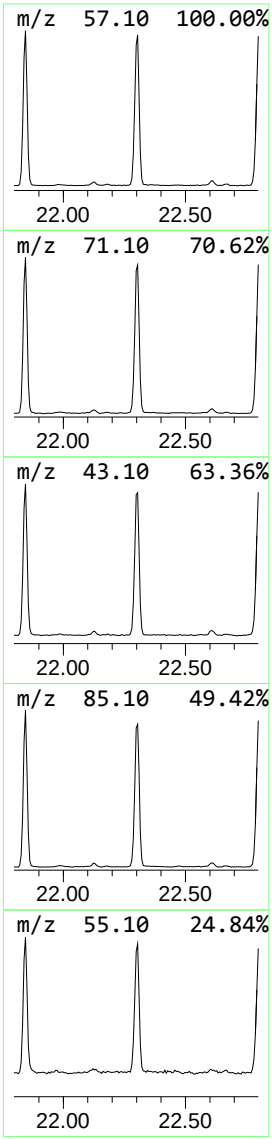
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.300	38.70 ng/ul	12336200	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricosane	324	C23H48	000638-67-5	96
2	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	94
3	Eicosane, 9-octyl-	394	C28H58	013475-77-9	94
4	Heptacosane	380	C27H56	000593-49-7	94
5	Hexacosane	366	C26H54	000630-01-3	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100720\
Data File : BN012096.D
Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

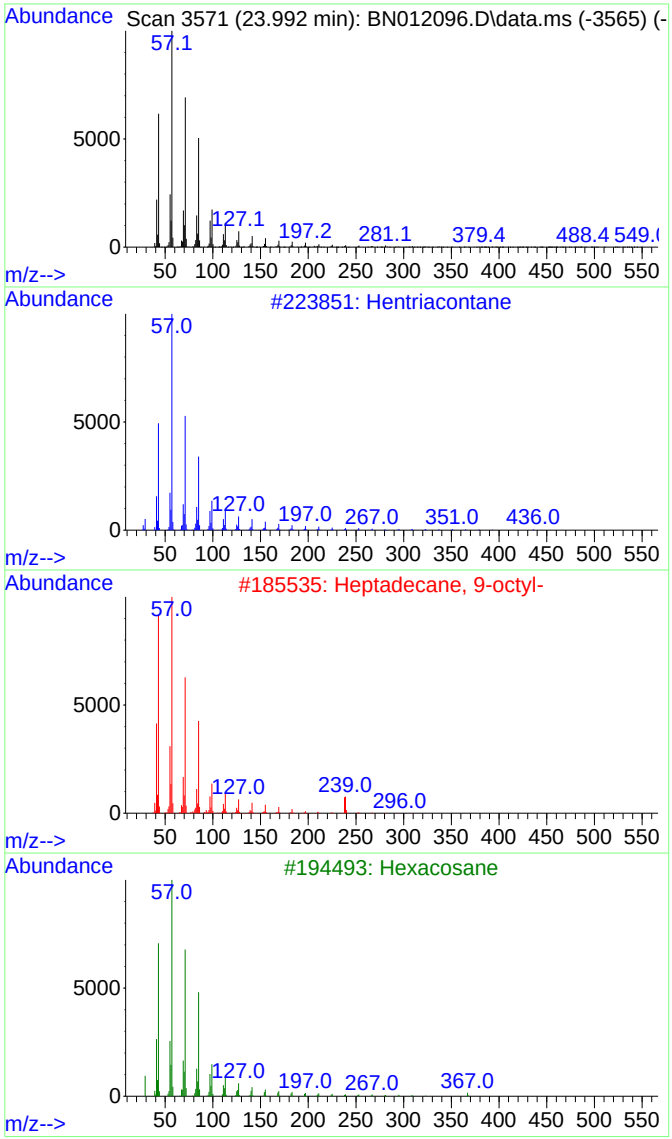
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

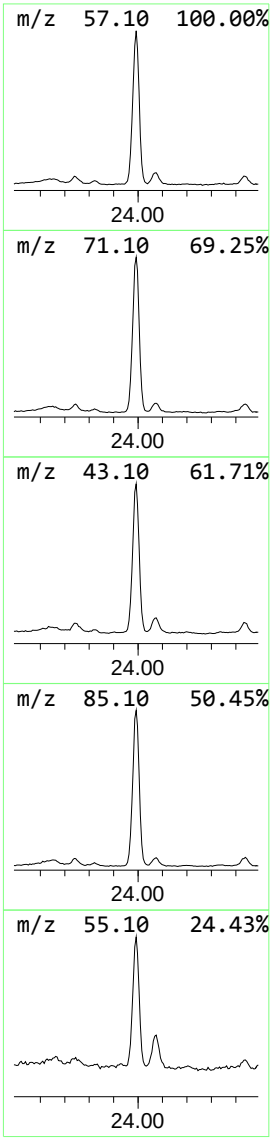
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.990	48.75 ng/ul	11600400	Perylene-d12	23.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hentriacontane	437	C31H64	000630-04-6	95
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
3	Hexacosane	366	C26H54	000630-01-3	91
4	2-methyloctacosane	408	C29H60	1000376-72-8	91
5	Heneicosane	296	C21H44	000629-94-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100720\
Data File : BN012096.D
Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AP4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

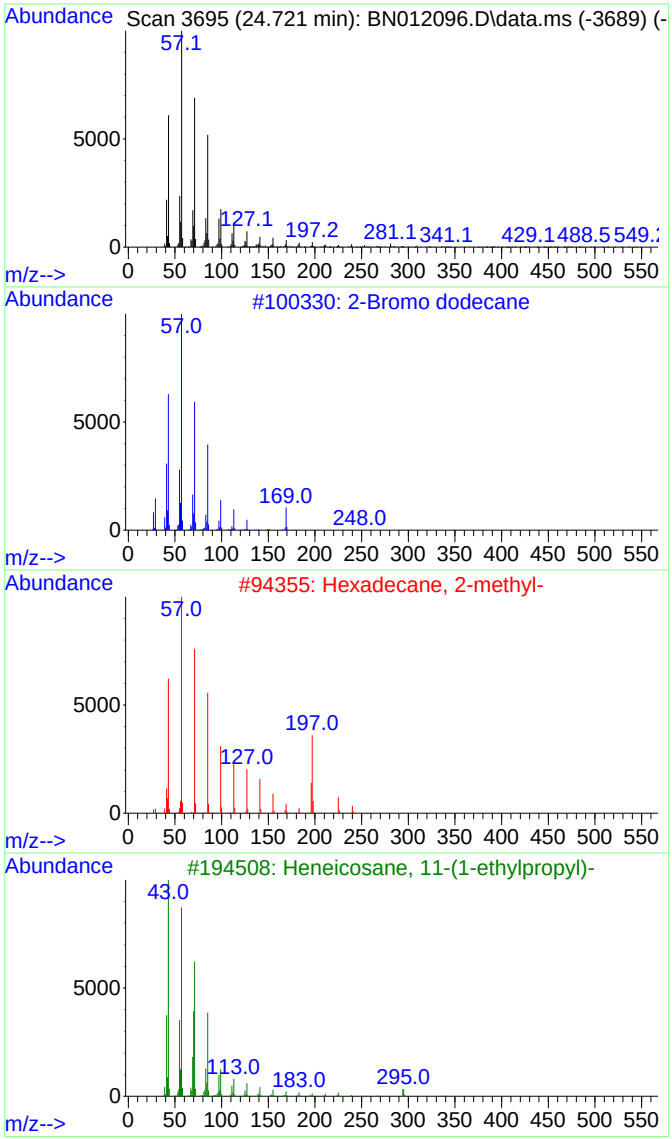
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

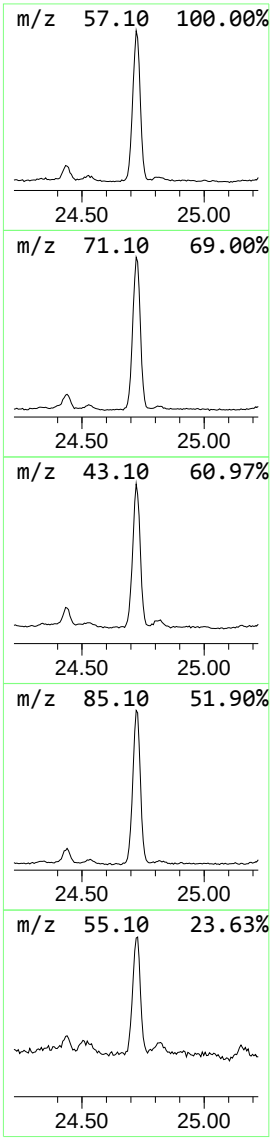
****

Peak Number 26 2-Bromo dodecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.720	32.64 ng/ul	7765390	Perylene-d12	23.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo dodecane	248	C12H25Br	013187-99-0	94
2	Hexadecane, 2-methyl-	240	C17H36	001560-92-5	93
3	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
4	2-methyloctacosane	408	C29H60	1000376-72-8	91
5	Octacosane	394	C28H58	000630-02-4	91





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Data File : BN012096.D
Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AP4

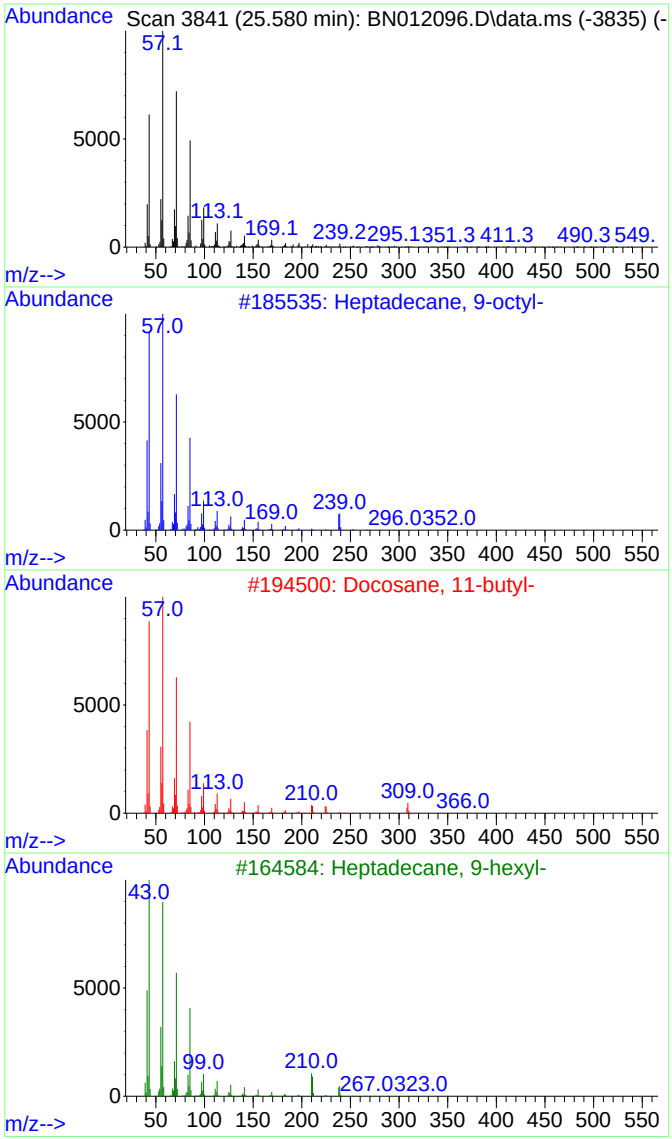
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

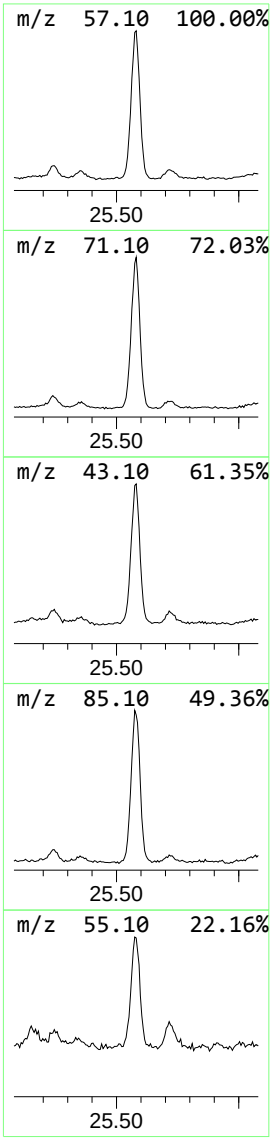
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.580	24.45 ng/ul	5817020	Perylene-d12	23.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2	Docosane, 11-butyl-	366	C26H54	013475-76-8	93
3	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	93
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100720\
Data File : BN012096.D
Acq On : 07 Oct 2020 14:35
Operator : CG/JU
Sample : L4184-20
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AP4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.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
Bicyclo[4.2.0]o...	5.670	116.5	ng/ul	13958300	1	7.652	2395630	20.0
Benzene, (1-met...	6.160	5.5	ng/ul	663106	1	7.652	2395630	20.0
Benzene, propyl-	6.650	2.6	ng/ul	313448	1	7.652	2395630	20.0
Ethanol, 2-(2-e...	7.270	2.0	ng/ul	246078	1	7.652	2395630	20.0
Tetramethylbuta...	7.730	2.1	ng/ul	248889	1	7.652	2395630	20.0
Benzene, 1,1'-(...	16.230	2.0	ng/ul	494091	4	17.045	4829030	20.0
[2.2]Paracyclop...	16.650	3.9	ng/ul	928537	4	17.045	4829030	20.0
n-Hexadecanoic ...	17.950	11.7	ng/ul	2818780	4	17.045	4829030	20.0
Fumaric acid, 2...	18.820	9.7	ng/ul	2351080	4	17.045	4829030	20.0
Octadecanoic acid	19.250	8.0	ng/ul	2559160	5	21.251	6376130	20.0
(DEL) Alkane: S...	20.010	2.0	ng/ul	649159	5	21.251	6376130	20.0
Octadecanoic ac...	20.450	2.2	ng/ul	700510	5	21.251	6376130	20.0
(DEL) Alkane: S...	20.510	6.0	ng/ul	1905790	5	21.251	6376130	20.0
Phosphoric acid...	20.730	4.6	ng/ul	1465960	5	21.251	6376130	20.0
(DEL) Alkane: S...	20.970	19.9	ng/ul	6343640	5	21.251	6376130	20.0
2-Bromomethyl-2...	21.030	2.9	ng/ul	934900	5	21.251	6376130	20.0
(DEL) Alkane: S...	21.420	31.7	ng/ul	10099500	5	21.251	6376130	20.0
(2,3-Diphenylcy...	21.470	13.6	ng/ul	4339500	5	21.251	6376130	20.0
1-Propene, 3-(2...	21.520	5.0	ng/ul	1611240	5	21.251	6376130	20.0
unknown-01	21.550	6.7	ng/ul	2138790	5	21.251	6376130	20.0
(DEL) Alkane: S...	21.850	35.3	ng/ul	11255000	5	21.251	6376130	20.0
(DEL) Alkane: S...	22.300	38.7	ng/ul	12336200	5	21.251	6376130	20.0
(DEL) Alkane: S...	23.990	48.8	ng/ul	11600400	6	23.463	4758890	20.0
2-Bromo dodecane	24.720	32.6	ng/ul	7765390	6	23.463	4758890	20.0
(DEL) Alkane: S...	25.580	24.4	ng/ul	5817020	6	23.463	4758890	20.0