LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015064.D
Acq On : 25 Jun 2021 06:57
Operator : CG/JU
Sample : M2720-01
Misc : GCMS CONFIRMATION
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0CP2

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: BN015064.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.134	4	8	11	rBV	68990	71933	1.60%	0.332%
2	3.163	11	13	15	rVV	28535	30692	0.68%	0.142%
3	3.187	15	17	19	rVV	46185	46635	1.04%	0.215%
4	3.222	19	23	37	rVB	4234395	4500554	100.00%	20.761%
5	3.516	69	73	74	rBV2	12958	13031	0.29%	0.060%
6	3.546	74	78	83	rVB	133573	154393	3.43%	0.712%
7	3.634	89	93	97	rVB2	9376	11489	0.26%	0.053%
8	3.911	134	140	146	rVB2	6334	9629	0.21%	0.044%
9	3.999	150	155	165	rVB	44586	55557	1.23%	0.256%
10	5.275	368	372	377	rBV3	4420	6970	0.15%	0.032%
11	6.369	554	558	564	rBV	9397	12435	0.28%	0.057%
12	7.722	781	788	796	rBV	839549	1291885	28.71%	5.959%
13	10.498	1253	1260	1267	rBV	1040836	1738354	38.63%	8.019%
14	14.351	1908	1915	1923	rBV	1456814	2068331	45.96%	9.541%
15	16.880	2341	2345	2350	rVB	6156	7680	0.17%	0.035%
16	17.098	2376	2382	2394	rBV	1613480	2234809	49.66%	10.309%
17	17.198	2394	2399	2402	rVV	4222	4810	0.11%	0.022%
18	17.275	2410	2412	2418	rBV6	4436	7038	0.16%	0.032%
19	17.698	2479	2484	2492	rVB3	27708	51739	1.15%	0.239%
20	17.822	2503	2505	2509	rVB3	5237	5591	0.12%	0.026%
21	17.951	2524	2527	2530	rBV3	10986	12608	0.28%	0.058%
22	18.016	2534	2538	2543	rVB2	3522	6355	0.14%	0.029%
23	18.174	2560	2565	2571	rBV	178153	228951	5.09%	1.056%
24	18.233	2571	2575	2580	rVV	42922	52993	1.18%	0.244%
25	18.274	2580	2582	2586	rVB3	10531	10538	0.23%	0.049%
26	18.457	2609	2613	2618	rVB	82699	103960	2.31%	0.480%
27	18.510	2618	2622	2625	rBV4	6604	7349	0.16%	0.034%
28	18.598	2633	2637	2640	rBV4	7612	9835	0.22%	0.045%
29	18.633	2640	2643	2647	rVB	28644	33943	0.75%	0.157%
30	18.733	2658	2660	2663	rVB4	5234	5670	0.13%	0.026%
31	18.945	2692	2696	2700	rBV2	36441	41080	0.91%	0.190%
32	18.998	2700	2705	2707	rBV	106287	145462	3.23%	0.671%
33	19.027	2707	2710	2716	rVV3	283879	379162	8.42%	1.749%
34	19.110	2720	2724	2728	rBV5	38936	47729	1.06%	0.220%
35	19.157	2728	2732	2738	rBV	9312	11519	0.26%	0.053%
36	19.233	2741	2745	2753	rBV4	62508	103730	2.30%	0.479%

LSC Area Percent Report

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Sample : M2720-01
Misc : GCMS CONFIRMATION
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0CP2

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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Peak separation: 5

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

37	19.310	2753	2758	2765	rVV	447585	581839	12.93%	2.684%
38	19.374	2765	2769	2773	rVB	155497	182039	4.04%	0.840%
39	19.445	2779	2781	2784	rVB4	5478	4953	0.11%	0.023%
40	19.504	2788	2791	2793	rBV2	14315	15355	0.34%	0.071%
41	19.574	2798	2803	2807	rVV	115290	142061	3.16%	0.655%
42	19.651	2809	2816	2820	rVV	194728	244783	5.44%	1.129%
43	19.704	2820	2825	2828	rVV	63491	76896	1.71%	0.355%
44	19.757	2828	2834	2840	rVB	442780	574280	12.76%	2.649%
45	19.904	2853	2859	2861	rBV3	44766	83516	1.86%	0.385%
46	19.927	2861	2863	2868	rVV4	25411	34321	0.76%	0.158%
47	19.974	2868	2871	2874	rVV2	16230	15624	0.35%	0.072%
48	20.021	2874	2879	2883	rVV2	193434	249014	5.53%	1.149%
49	20.074	2883	2888	2893	rVB	386655	456669	10.15%	2.107%
50	20.151	2898	2901	2904	rBV3	17823	17721	0.39%	0.082%
51	20.227	2908	2914	2917	rBV6	34334	52489	1.17%	0.242%
52	20.304	2922	2927	2931	rVV	223908	262539	5.83%	1.211%
53	20.357	2931	2936	2938	rVV	113592	155710	3.46%	0.718%
54	20.380	2938	2940	2945	rVB	172188	177383	3.94%	0.818%
55	20.457	2950	2953	2957	rVB3	45979	53196	1.18%	0.245%
56	20.521	2961	2964	2967	rBV3	15768	20476	0.45%	0.094%
57	20.621	2973	2981	2987	rBV	273266	456837	10.15%	2.107%
58	20.704	2992	2995	2999	rBV6	16307	17617	0.39%	0.081%
59	20.768	3004	3006	3010	rVB3	9963	10168	0.23%	0.047%
60	20.927	3029	3033	3037	rVB2	86210	104728	2.33%	0.483%
61	21.039	3049	3052	3055	rVB5	14104	14850	0.33%	0.069%
62	21.174	3072	3075	3079	rBV3	39215	38761	0.86%	0.179%
63	21.292	3089	3095	3100	rBV	1652654	2068485	45.96%	9.542%
64	21.633	3151	3153	3158	rVB6	21894	27150	0.60%	0.125%
65	22.880	3362	3365	3370	rVB	46456	62594	1.39%	0.289%
66	23.539	3470	3477	3486	rBV2	985999	1877514	41.72%	8.661%
67	24.104	3569	3573	3579	rVB3	40900	73934	1.64%	0.341%

Sum of corrected areas: 21677941

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ALS Vial : 28 Sample Multiplier: 1

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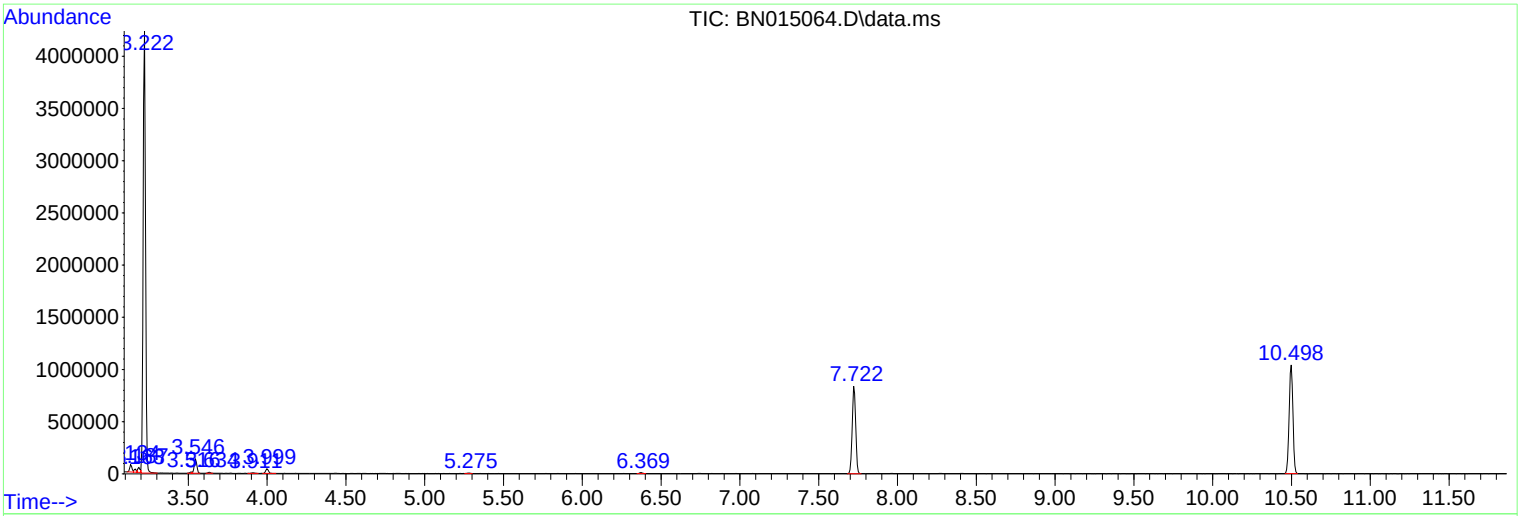
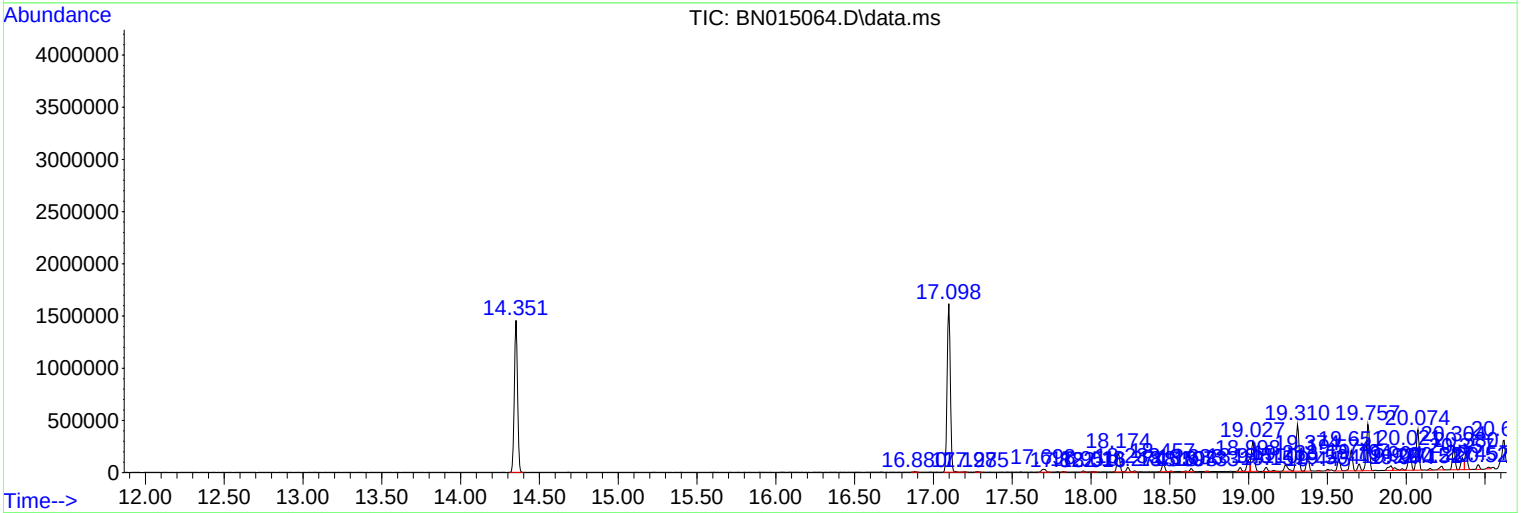
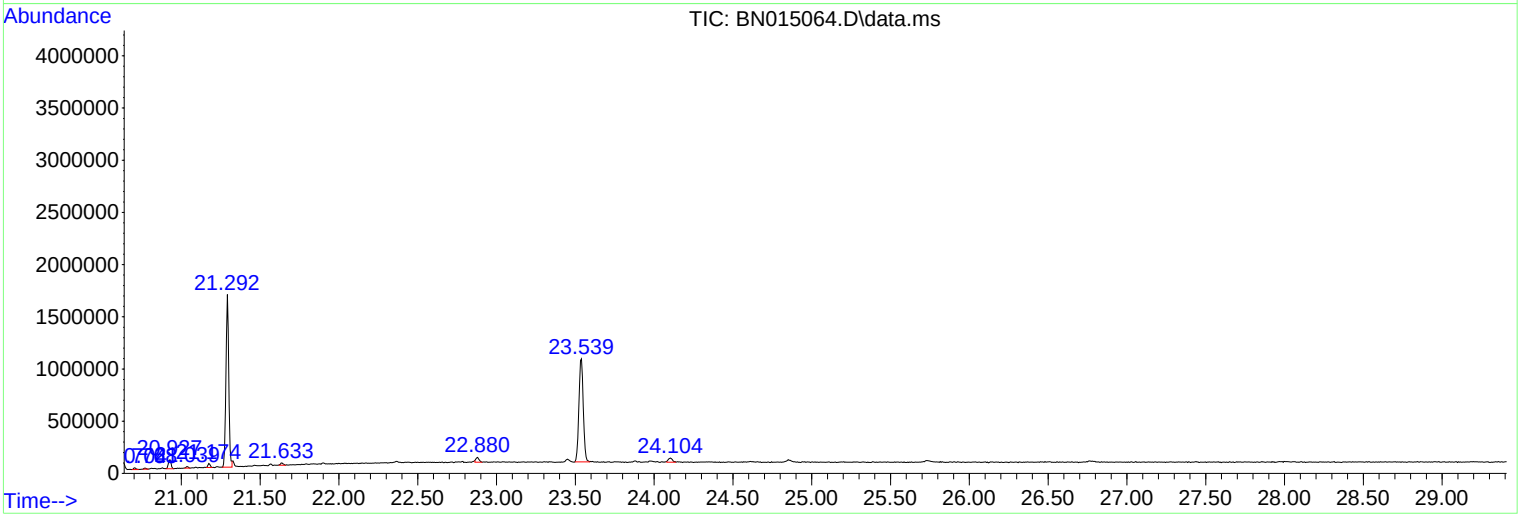
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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







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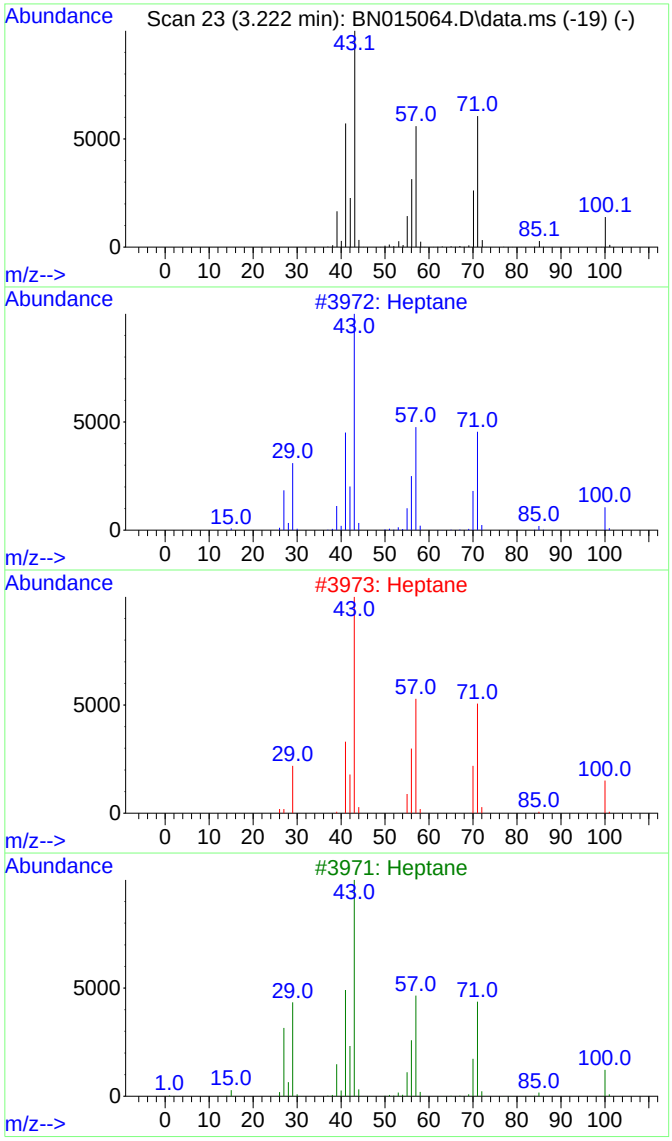
Quant Title : SVOA CALIBRATION

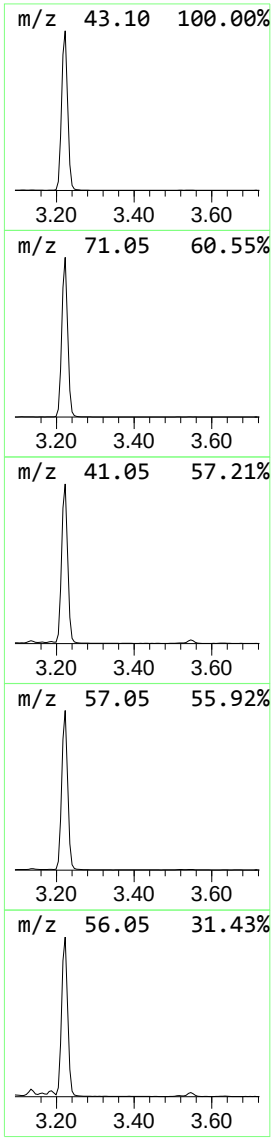
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.222	69.67 ng/ul	4500550	1,4-Dichlorobenzene-d4	7.722

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane	100	C7H16	000142-82-5	95
2	Heptane	100	C7H16	000142-82-5	91
3	Heptane	100	C7H16	000142-82-5	90
4	Heptane	100	C7H16	000142-82-5	86
5	Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	59





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ALS Vial : 28 Sample Multiplier: 1

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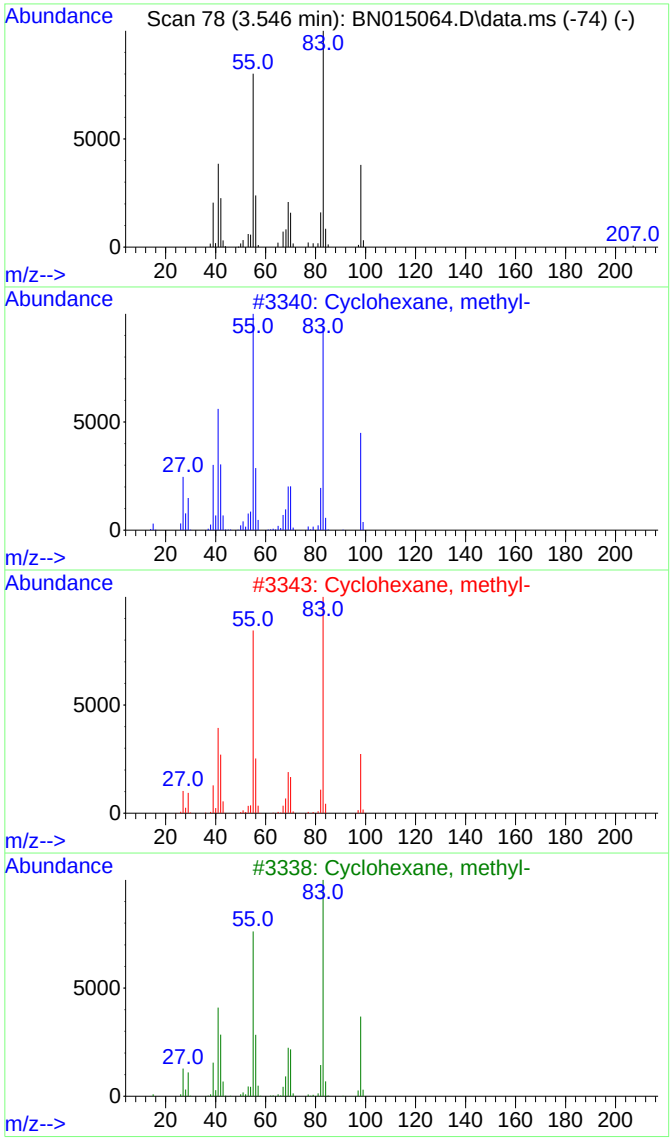
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

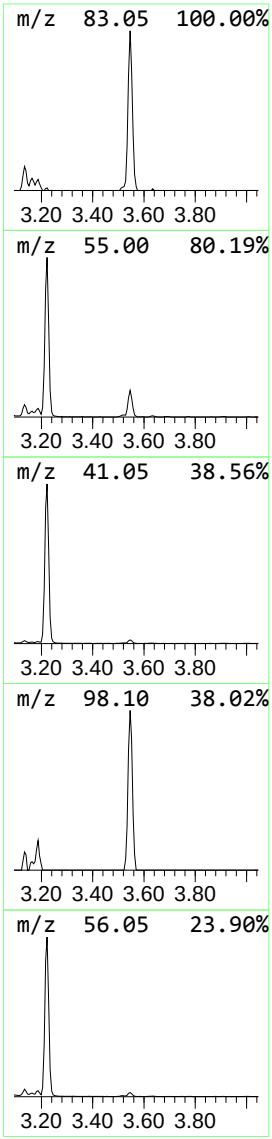
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Cyclic3.546 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.546	2.39 ng/ul	154393	1,4-Dichlorobenzene-d4	7.722

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, methyl-	98	C7H14	000108-87-2	93
2	Cyclohexane, methyl-	98	C7H14	000108-87-2	91
3	Cyclohexane, methyl-	98	C7H14	000108-87-2	91
4	Cyclohexane, methyl-	98	C7H14	000108-87-2	87
5	2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	58





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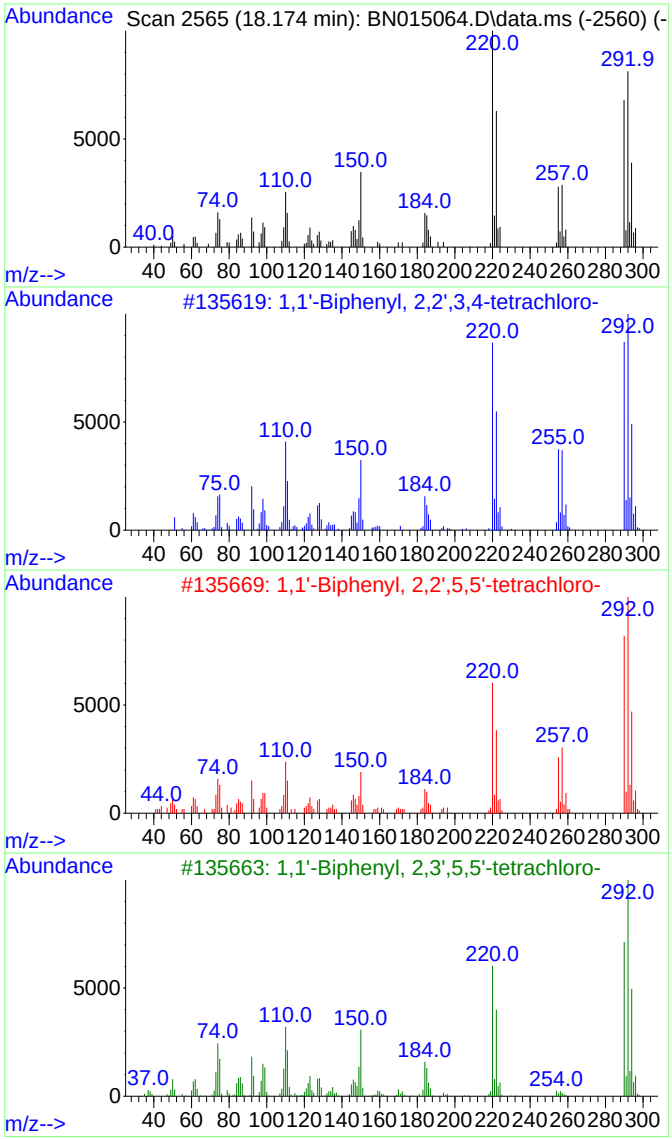
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

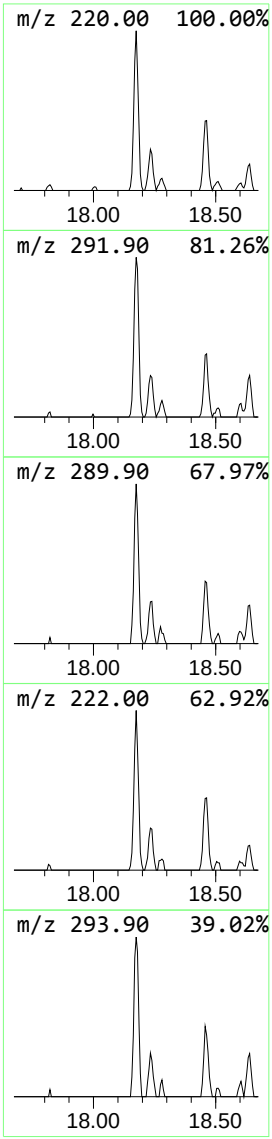
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1,1'-Biphenyl, 2,2',3,4-tet... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.174	2.05 ng/ul	228951	Phenanthrene-d10	17.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	99
2	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
3	1,1'-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	99
4	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
5	2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	99





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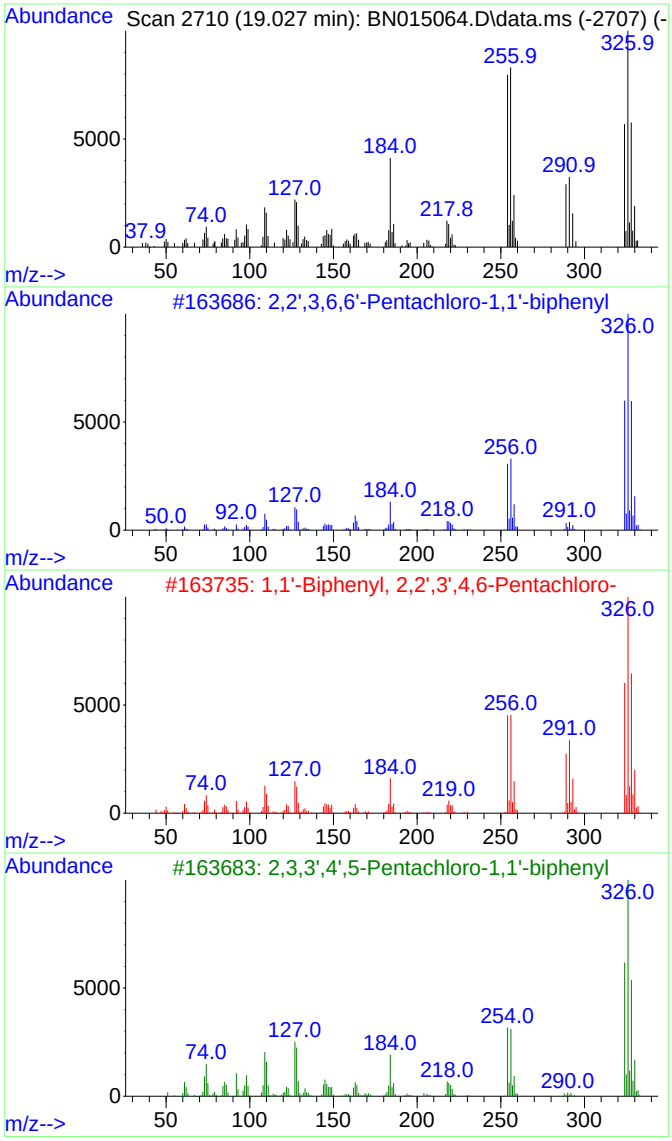
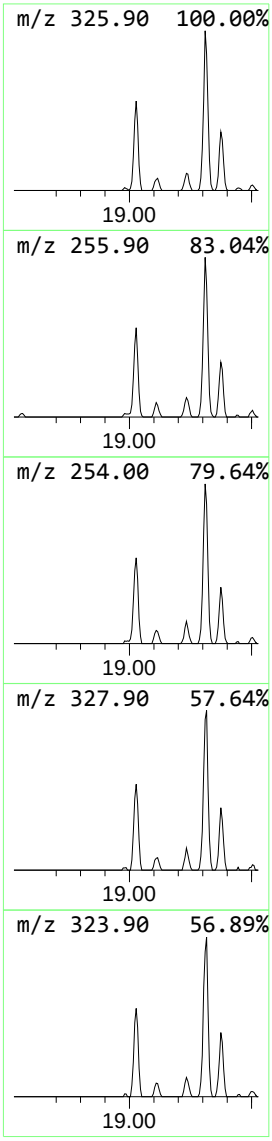
****
Peak Number 4 2,2',3,6,6'-Pentachloro-1,1... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.027	3.39 ng/ul	379162	Phenanthrene-d10	17.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,6,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-54-9	99
2	1,1'-Biphenyl, 2,2',3',4,6-Penta...	324	C12H5Cl5	060233-25-2	97
3	2,3,3',4',5-Pentachloro-1,1'-bip...	324	C12H5Cl5	070424-68-9	97
4	1,1'-Biphenyl, 2,2',4,5,6'-penta...	324	C12H5Cl5	068194-06-9	96
5	1,1'-Biphenyl, 2,2',3,3',4-penta...	324	C12H5Cl5	052663-62-4	96





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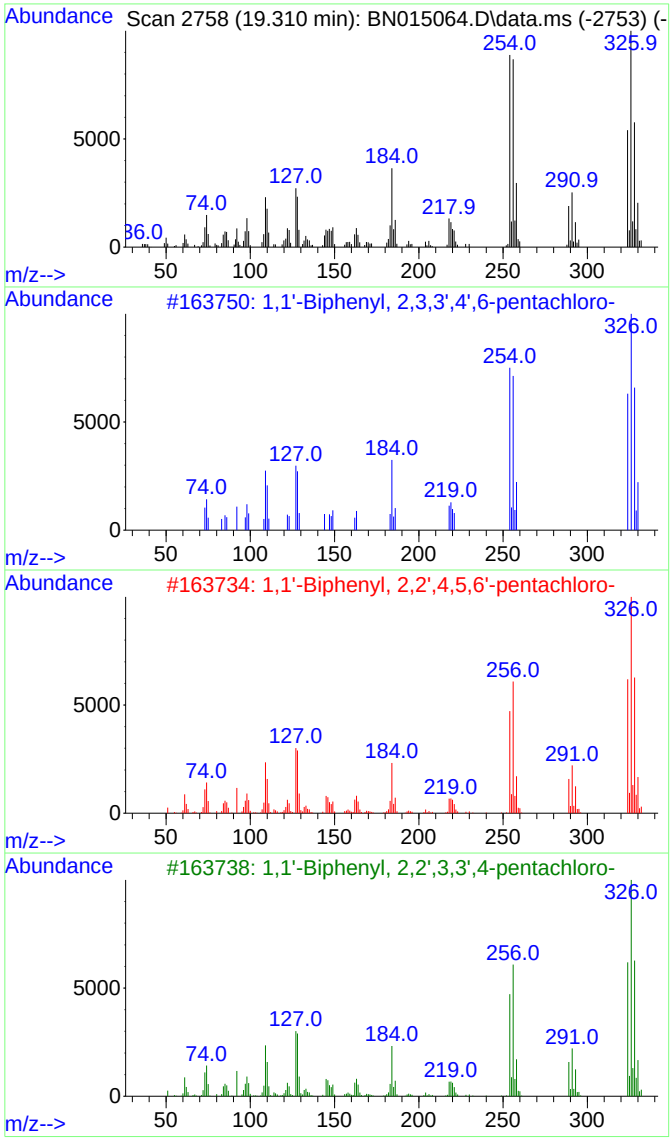
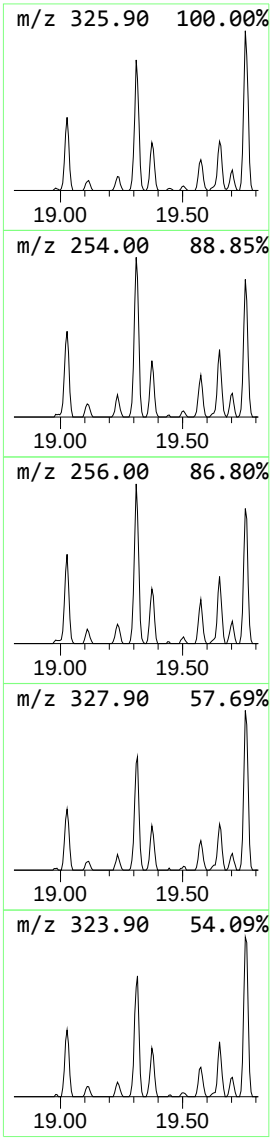
****
Peak Number 5 1,1'-Biphenyl, 2,3,3',4',6-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.310	5.63 ng/ul	581839	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99
2	1,1'-Biphenyl, 2,2',4,5,6'-penta...	324	C12H5Cl5	068194-06-9	99
3	1,1'-Biphenyl, 2,2',3,3',4-penta...	324	C12H5Cl5	052663-62-4	99
4	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99
5	1,1'-Biphenyl, 2,2',4,6,6'-Penta...	324	C12H5Cl5	056558-16-8	99





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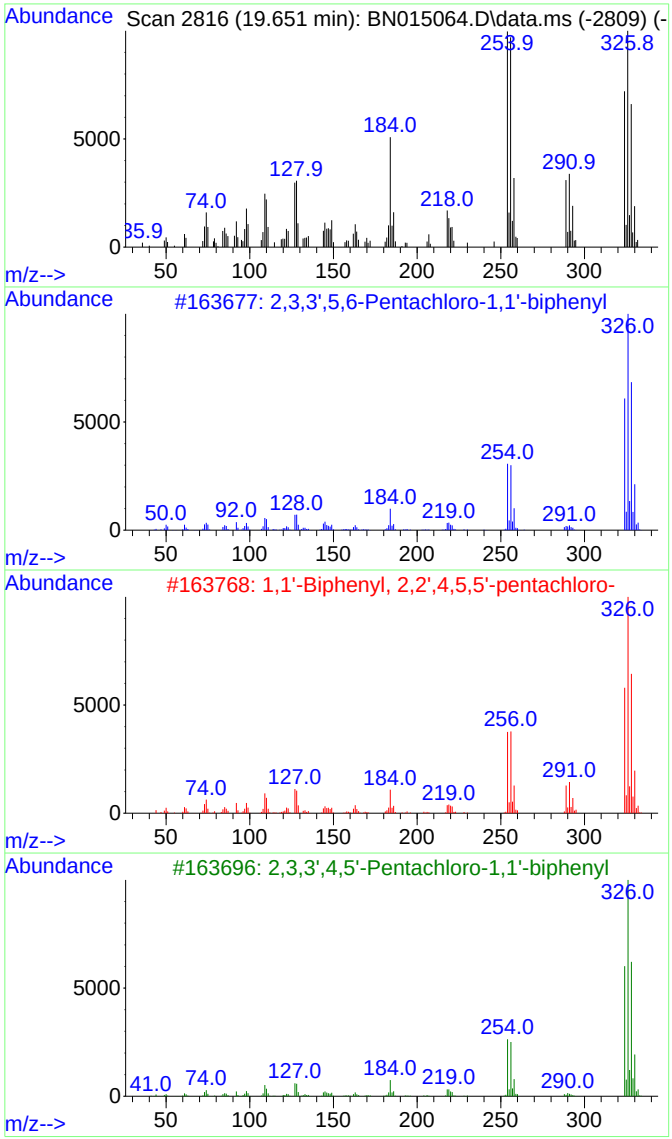
Quant Title : SVOA CALIBRATION

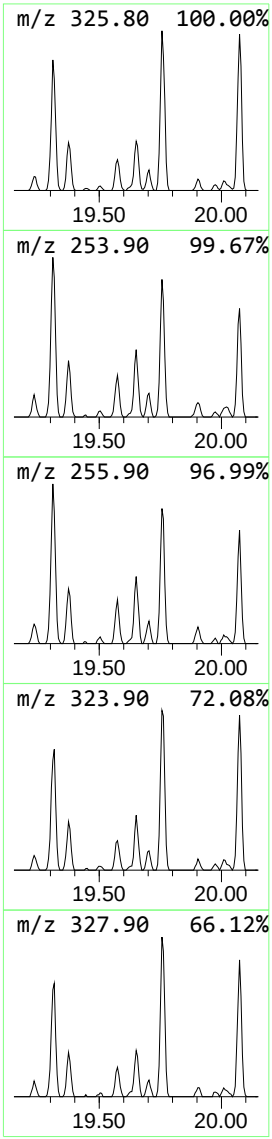
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 2,3,3',5,6-Pentachloro-1,1'... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.651	2.37 ng/ul	244783	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	99
2	1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	99
3	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	99
4	2,2',3,4,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-57-2	99
5	1,1'-Biphenyl, 2,2',4,5,6'-penta...	324	C12H5Cl5	068194-06-9	99





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Sample : M2720-01
Misc : GCMS CONFIRMATION
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0CP2

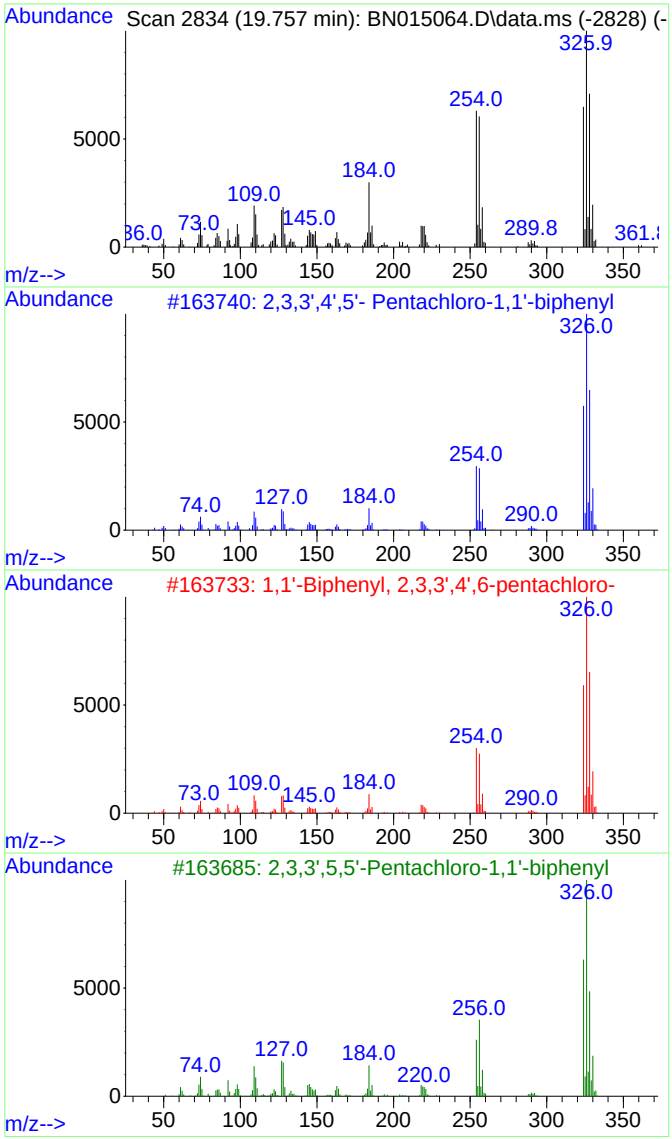
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

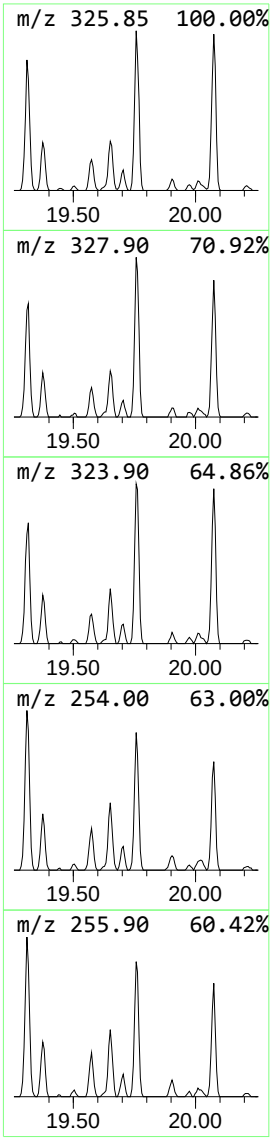
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 2,3,3',4',5'- Pentachloro-1... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.757	5.55 ng/ul	574280	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,3',4',5'- Pentachloro-1,1'-b...	324	C12H5Cl5	076842-07-4	99
2	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99
3	2,3,3',5,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-32-0	99
4	1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99
5	3,3',4,5,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-33-1	99





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Data File : BN015064.D
Acq On : 25 Jun 2021 06:57
Operator : CG/JU
Sample : M2720-01
Misc : GCMS CONFIRMATION
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0CP2

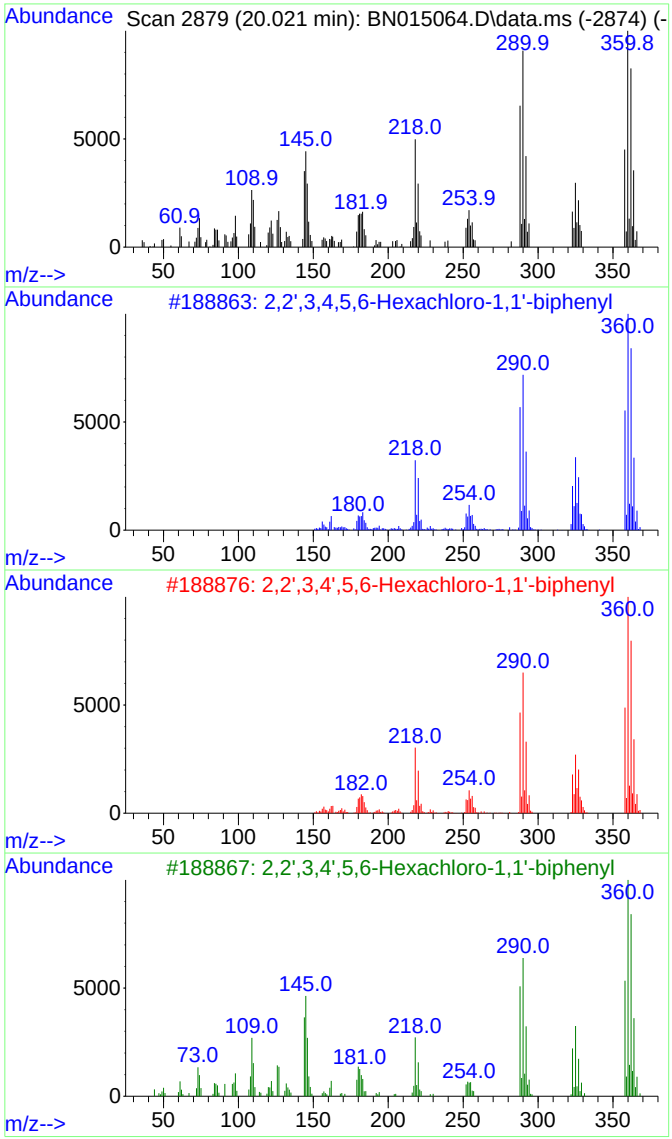
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

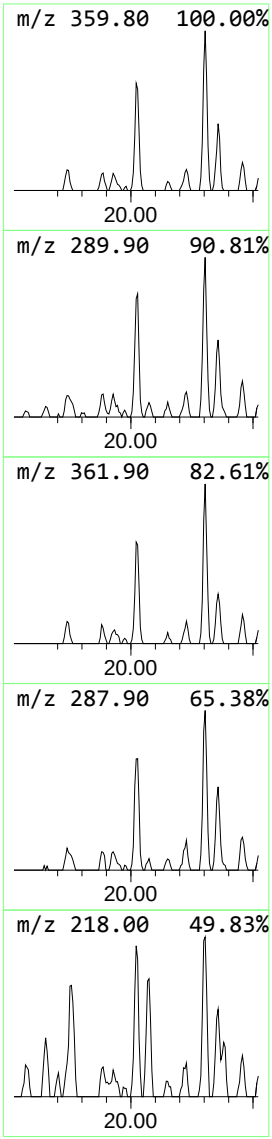
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 2,2',3,4,5,6-Hexachloro-1,1... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.021	2.41 ng/ul	249014	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4,5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-61-4	99
2	2,2',3,4',5,6-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-13-8	99
3	2,2',3,4',5,6-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-13-8	99
4	1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	99
5	2,3,3',4,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-43-8	99





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Data File : BN015064.D
Acq On : 25 Jun 2021 06:57
Operator : CG/JU
Sample : M2720-01
Misc : GCMS CONFIRMATION
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0CP2

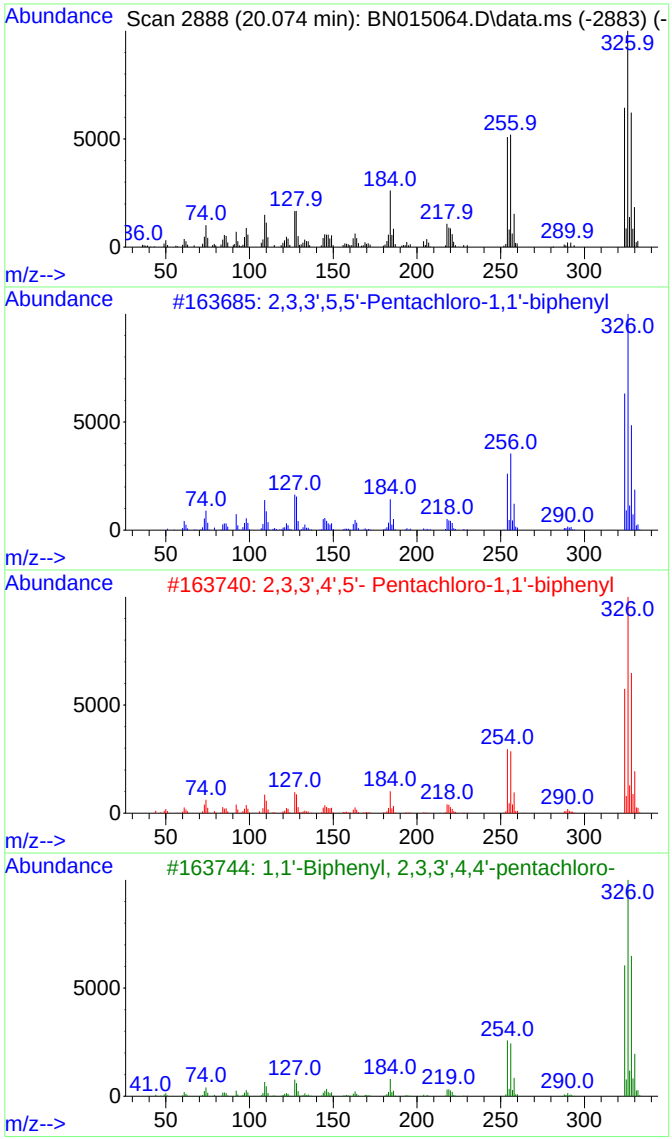
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

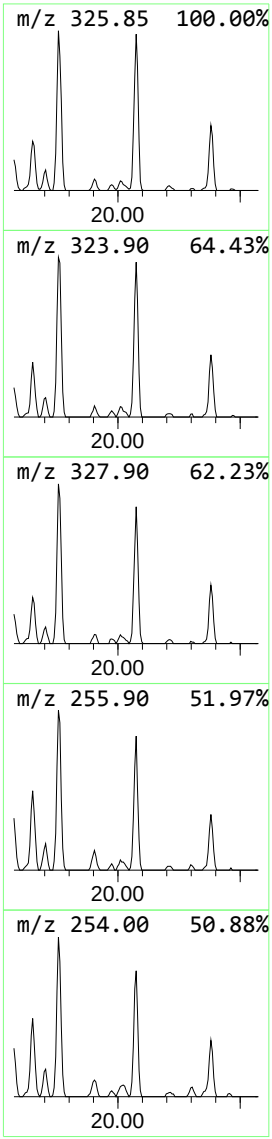
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 2,3,3',5,5'-Pentachloro-1,1... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.074	4.42 ng/ul	456669	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,3',5,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-32-0	99
2	2,3,3',4',5'- Pentachloro-1,1'-b...	324	C12H5Cl5	076842-07-4	99
3	1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99
4	3,3',4,4,',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	99
5	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99





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Data File : BN015064.D
Acq On : 25 Jun 2021 06:57
Operator : CG/JU
Sample : M2720-01
Misc : GCMS CONFIRMATION
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0CP2

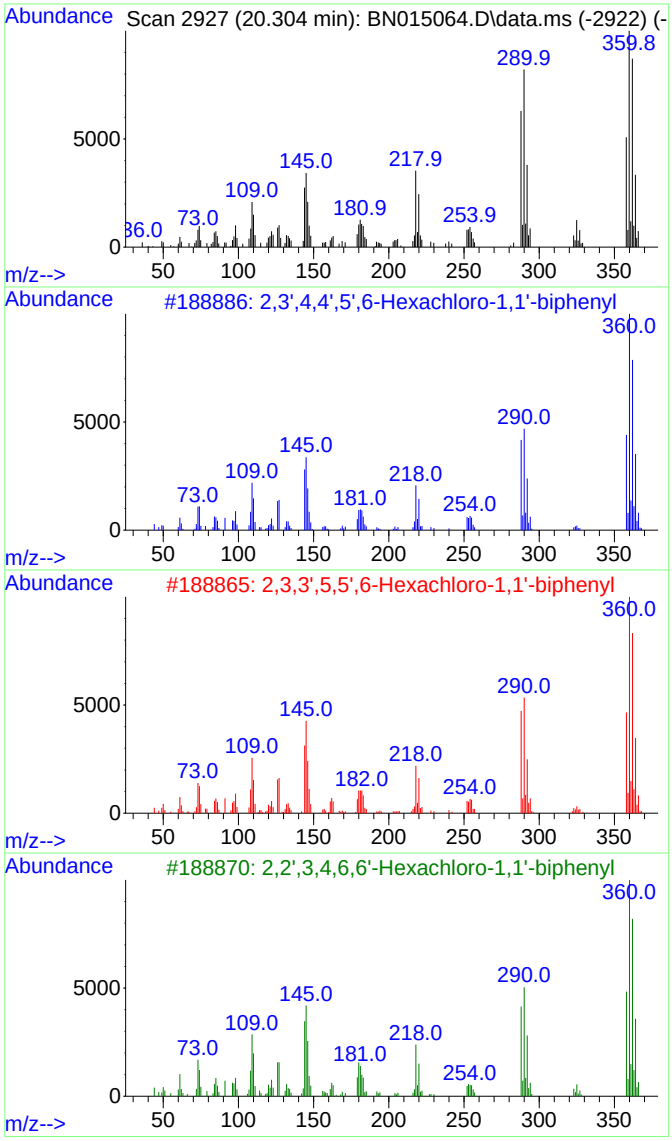
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

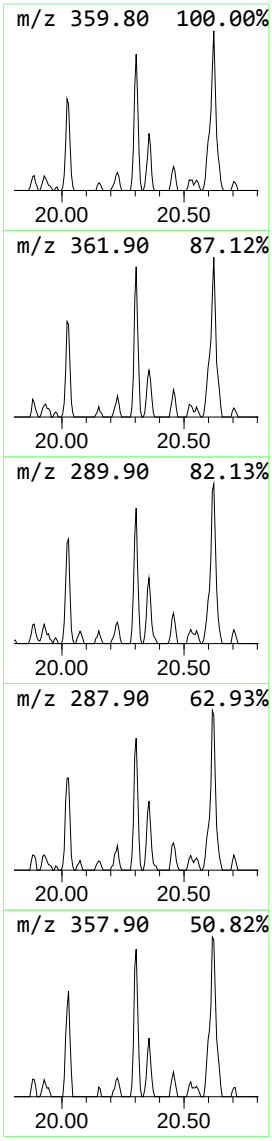
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 2,3',4,4',5',6-Hexachloro-1... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.304	2.54 ng/ul	262539	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3',4,4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	99
2	2,3,3',5,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-46-1	99
3	2,2',3,4,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-40-5	99
4	2,3,3',4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-45-0	99
5	2,3,3',4',5,5'-Hexachloro-1,1'-b...	358	C12H4Cl6	039635-34-2	99





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
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Sample : M2720-01
Misc : GCMS CONFIRMATION
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Instrument :
BNA_N
ClientSampleId :
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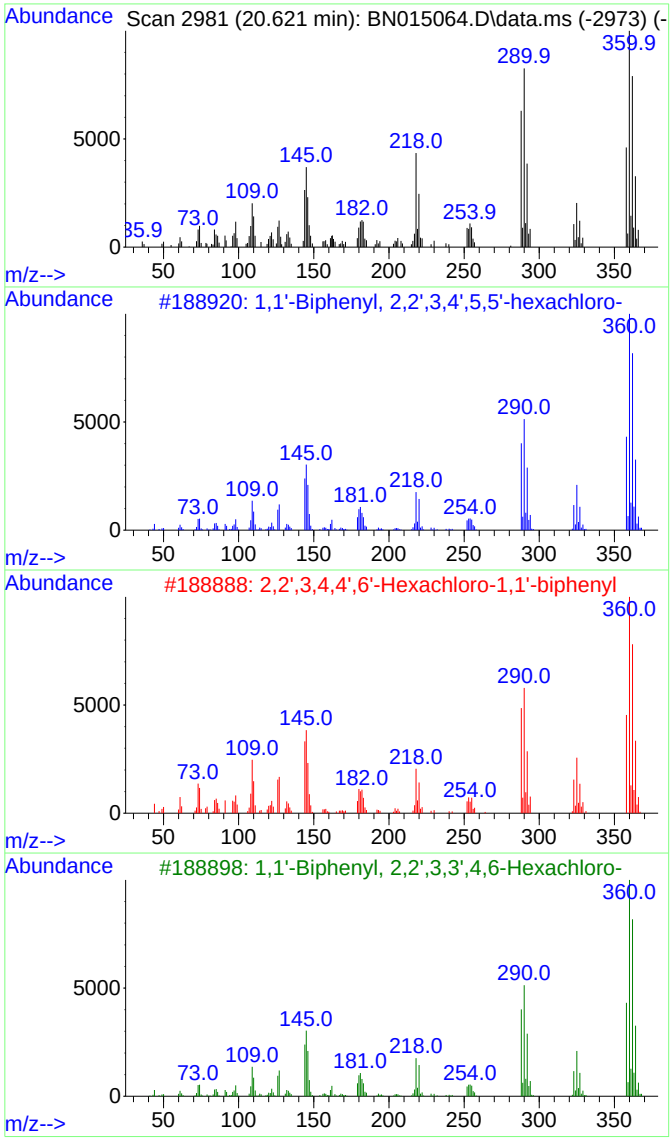
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

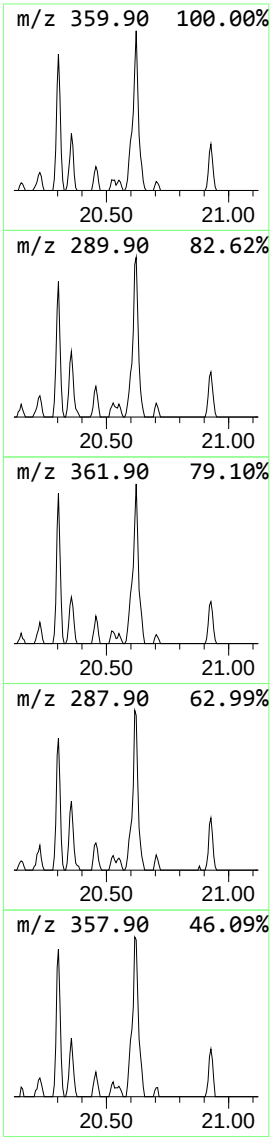
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 1,1'-Biphenyl, 2,2',3,4',5,... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.621	4.42 ng/ul	456837	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,4',5,5'-he...	358	C12H4Cl6	051908-16-8	99
2	2,2',3,4,4',6'-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-64-4	99
3	1,1'-Biphenyl, 2,2',3,3',4,6-Hex...	358	C12H4Cl6	061798-70-7	99
4	1,1'-Biphenyl, 2,2',4,4',5',6-He...	358	C12H4Cl6	060145-22-4	99
5	1,1'-Biphenyl, 2,3,3',4,4',5'-he...	358	C12H4Cl6	069782-90-7	99





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015064.D
Acq On : 25 Jun 2021 06:57
Operator : CG/JU
Sample : M2720-01
Misc : GCMS CONFIRMATION
ALS Vial : 28 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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: S...	3.222	69.7	ng/ul	4500550	1	7.722	1291890	20.0
(DEL) Alkane: C...	3.546	2.4	ng/ul	154393	1	7.722	1291890	20.0
1,1'-Biphenyl, ...	18.174	2.0	ng/ul	228951	4	17.098	2234810	20.0
2,2',3,6,6'-Pen...	19.027	3.4	ng/ul	379162	4	17.098	2234810	20.0
1,1'-Biphenyl, ...	19.310	5.6	ng/ul	581839	5	21.292	2068490	20.0
2,3,3',5,6-Pent...	19.651	2.4	ng/ul	244783	5	21.292	2068490	20.0
2,3,3',4',5'- P...	19.757	5.5	ng/ul	574280	5	21.292	2068490	20.0
2,2',3,4,5,6-He...	20.021	2.4	ng/ul	249014	5	21.292	2068490	20.0
2,3,3',5,5'-Pen...	20.074	4.4	ng/ul	456669	5	21.292	2068490	20.0
2,3',4,4',5',6-...	20.304	2.5	ng/ul	262539	5	21.292	2068490	20.0
1,1'-Biphenyl, ...	20.621	4.4	ng/ul	456837	5	21.292	2068490	20.0