

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021722\
 Data File : BN018639.D
 Acq On : 17 Feb 2022 22:04
 Operator : CG/JU
 Sample : N1507-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGZF5

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-BN021622.M
 Title : SVOA CALIBRATION

Signal : TIC: BN018639.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.040	6	9	17	rVB	218951	247117	6.76%	1.271%
2	3.128	20	24	26	rVV	59379	68701	1.88%	0.353%
3	3.152	26	28	30	rVV	29155	29565	0.81%	0.152%
4	3.181	30	33	35	rVV	42978	44040	1.21%	0.226%
5	3.223	35	40	45	rVB	3549498	3652988	100.00%	18.783%
6	3.540	90	94	95	rBV3	9886	11315	0.31%	0.058%
7	3.570	95	99	105	rVB	108526	129069	3.53%	0.664%
8	3.664	112	115	119	rVB2	8957	10923	0.30%	0.056%
9	4.052	178	181	188	rVB	16094	23207	0.64%	0.119%
10	4.522	257	261	265	rVB3	4136	6757	0.18%	0.035%
11	4.728	293	296	301	rBV5	2371	4095	0.11%	0.021%
12	7.916	831	838	845	rBV	822988	1303327	35.68%	6.701%
13	10.410	1258	1262	1270	rVB5	11252	16731	0.46%	0.086%
14	10.722	1307	1315	1328	rBV	1047872	1746313	47.81%	8.979%
15	10.999	1358	1362	1365	rVB5	2811	4177	0.11%	0.021%
16	11.816	1497	1501	1506	rBV4	4118	6125	0.17%	0.031%
17	14.345	1927	1931	1935	rBV5	2410	3718	0.10%	0.019%
18	14.551	1958	1966	1975	rBV	1346648	2033841	55.68%	10.457%
19	16.287	2257	2261	2265	rVB3	3520	5206	0.14%	0.027%
20	17.075	2390	2395	2398	rBV2	5039	7501	0.21%	0.039%
21	17.286	2424	2431	2442	rBV	1417209	2039612	55.83%	10.487%
22	17.381	2445	2447	2451	rVB2	5347	7176	0.20%	0.037%
23	17.898	2533	2535	2541	rVB5	6347	7534	0.21%	0.039%
24	17.963	2543	2546	2549	rVB4	5656	6021	0.16%	0.031%
25	18.363	2608	2614	2619	rBV	141341	193276	5.29%	0.994%
26	18.422	2619	2624	2628	rVV	27833	36310	0.99%	0.187%
27	18.463	2628	2631	2637	rVB7	5483	8102	0.22%	0.042%
28	18.645	2658	2662	2666	rBV	51104	66278	1.81%	0.341%
29	18.822	2688	2692	2697	rVB6	14366	20348	0.56%	0.105%
30	19.128	2738	2744	2747	rBV7	14559	23975	0.66%	0.123%
31	19.175	2747	2752	2754	rVV	106299	142826	3.91%	0.734%
32	19.204	2754	2757	2763	rVV	299727	387578	10.61%	1.993%
33	19.286	2765	2771	2775	rVV5	41624	74347	2.04%	0.382%
34	19.339	2777	2780	2785	rVB	12349	18077	0.49%	0.093%
35	19.410	2788	2792	2800	rBV4	67123	96215	2.63%	0.495%
36	19.486	2800	2805	2812	rVV	445380	581038	15.91%	2.988%

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Stop Thrs : 0

Peak Location: TOP

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37	19.551	2812	2816	2821	rVV	125669	155853	4.27%	0.801%
38	19.680	2834	2838	2840	rBV4	18775	24667	0.68%	0.127%
39	19.751	2846	2850	2855	rVB	103251	133557	3.66%	0.687%
40	19.828	2858	2863	2867	rVV	178480	219140	6.00%	1.127%
41	19.875	2867	2871	2874	rVV2	49080	61962	1.70%	0.319%
42	19.933	2874	2881	2886	rVB	380996	516621	14.14%	2.656%
43	20.057	2898	2902	2904	rBV4	37403	53936	1.48%	0.277%
44	20.145	2915	2917	2920	rVB4	15828	16685	0.46%	0.086%
45	20.198	2920	2926	2930	rBV2	179913	250079	6.85%	1.286%
46	20.245	2930	2934	2939	rVB	298716	352832	9.66%	1.814%
47	20.398	2956	2960	2962	rBV3	30945	38228	1.05%	0.197%
48	20.475	2968	2973	2977	rBV	200548	245478	6.72%	1.262%
49	20.527	2978	2982	2984	rVV	113883	149094	4.08%	0.767%
50	20.551	2984	2986	2992	rVV	107720	120478	3.30%	0.619%
51	20.627	2995	2999	3004	rBV3	49292	70554	1.93%	0.363%
52	20.792	3019	3027	3034	rVB	231922	402325	11.01%	2.069%
53	21.092	3075	3078	3084	rVB4	83458	111751	3.06%	0.575%
54	21.463	3136	3141	3146	rBV	1430811	1802026	49.33%	9.266%
55	23.863	3542	3549	3558	rVB2	799692	1660068	45.44%	8.536%

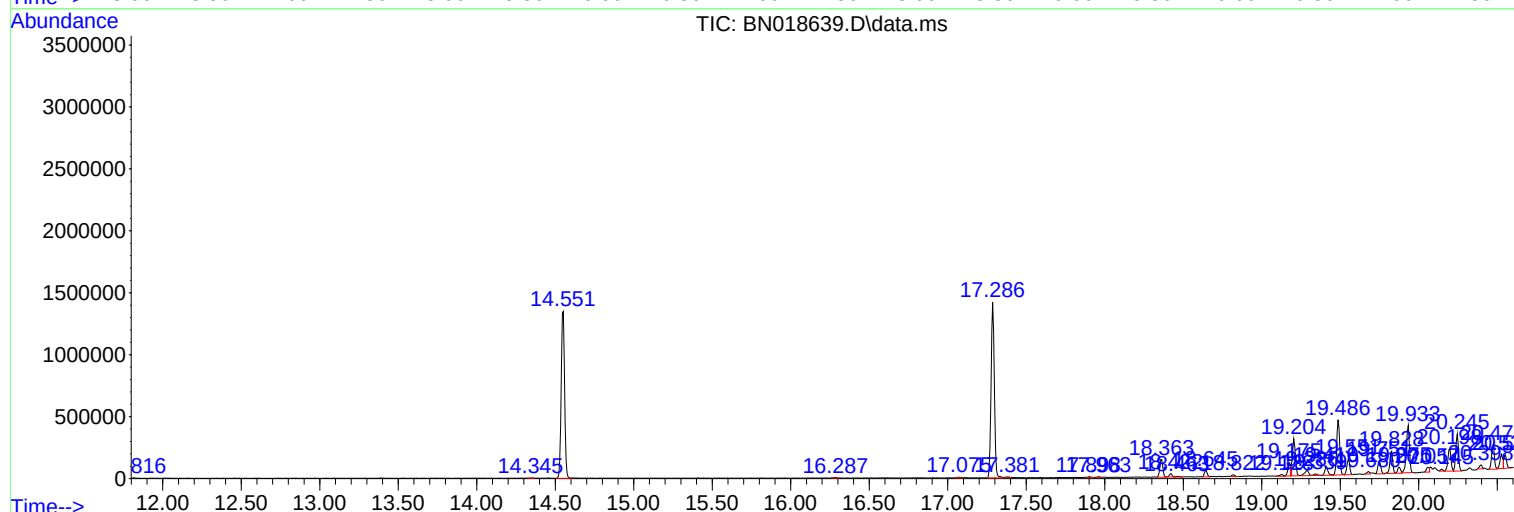
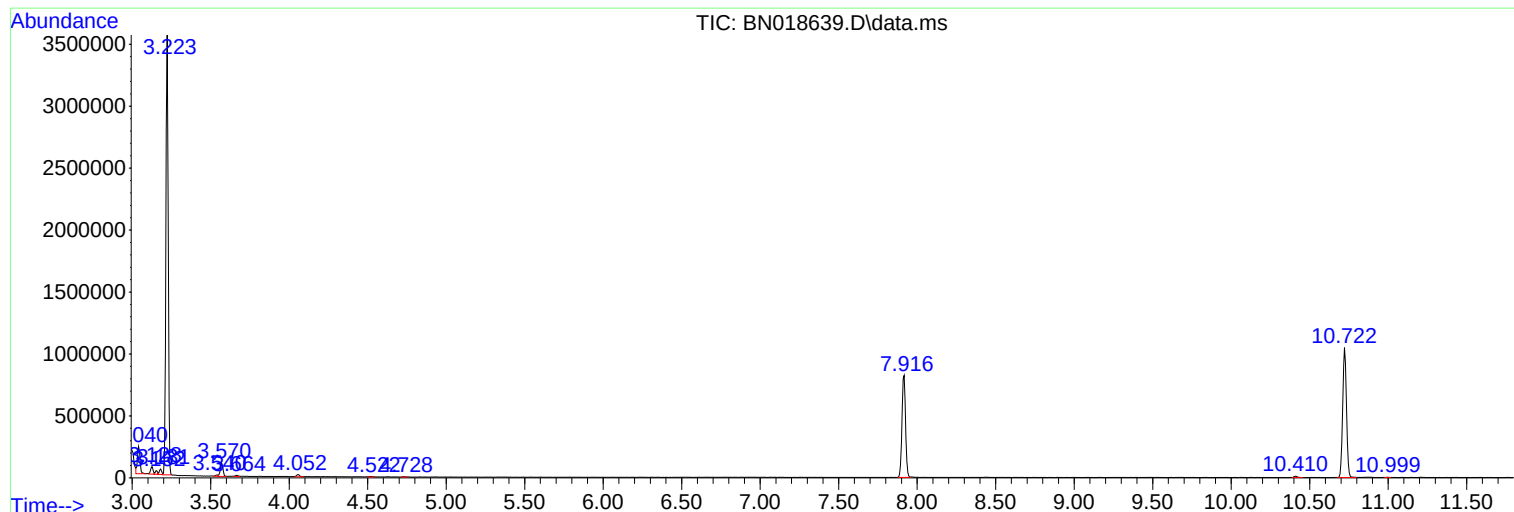
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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



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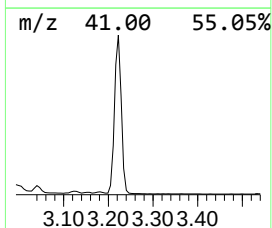
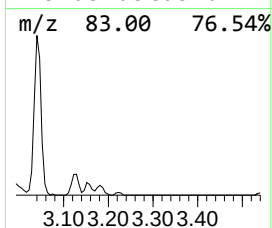
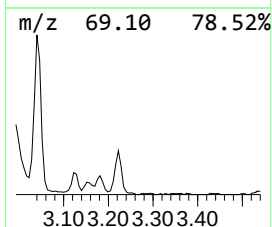
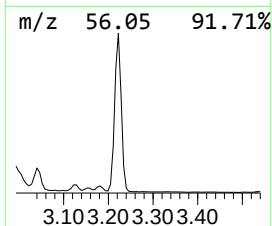
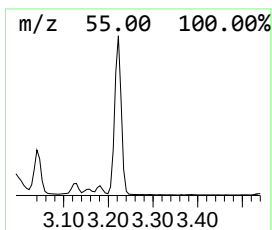
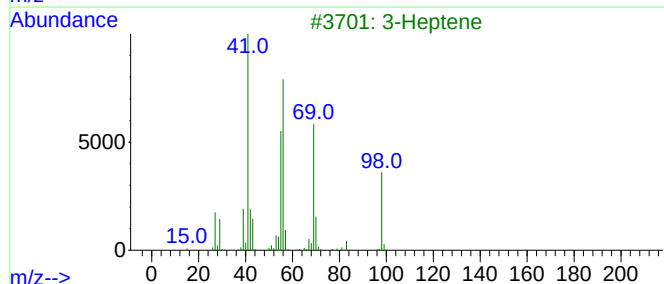
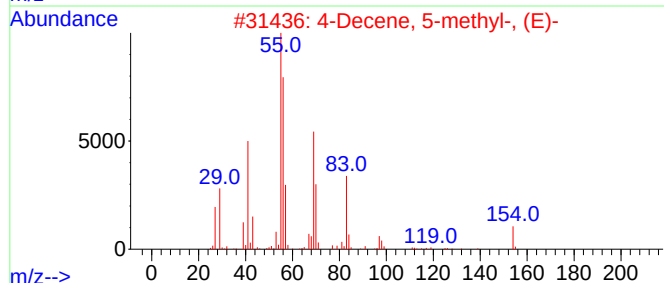
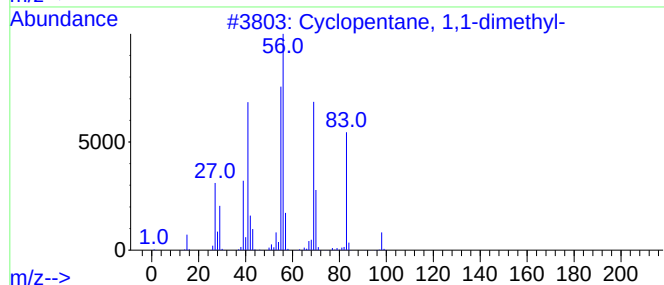
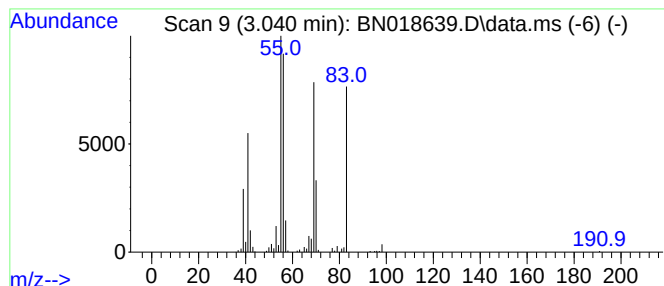
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.040 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.040	3.79 ng/ul	247117	1,4-Dichlorobenzene-d4	7.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	83
2			4-Decene, 5-methyl-, (E)-	154	C11H22	062338-51-6	59
3			3-Heptene	98	C7H14	000592-78-9	52
4			3-Undecene, 8-methyl-	168	C12H24	1000061-84-4	50
5			(R)-(+)-3-Methylcyclopentanone	98	C6H10O	006672-30-6	47



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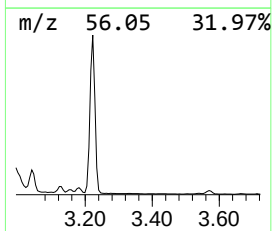
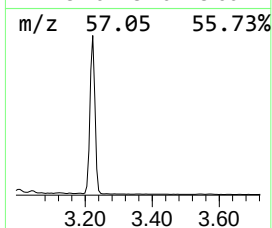
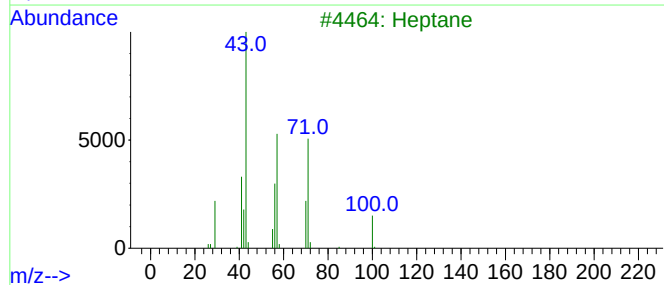
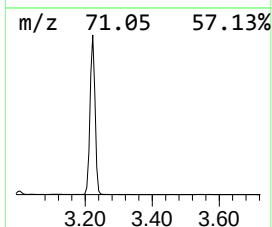
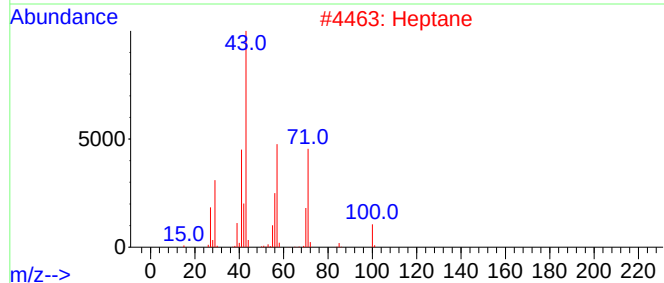
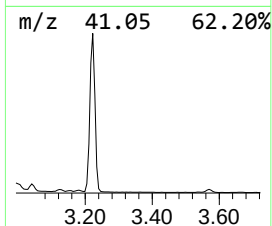
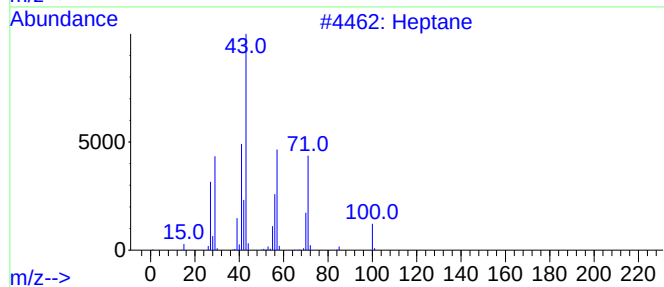
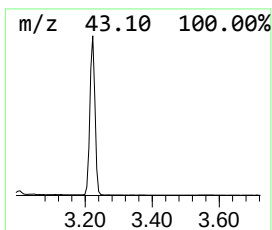
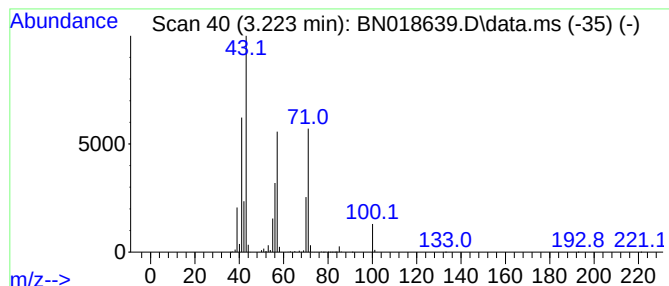
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.223	56.06 ng/ul	3652990	1,4-Dichlorobenzene-d4	7.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	94
2			Heptane	100	C7H16	000142-82-5	91
3			Heptane	100	C7H16	000142-82-5	91
4			Heptane	100	C7H16	000142-82-5	87
5			Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	45



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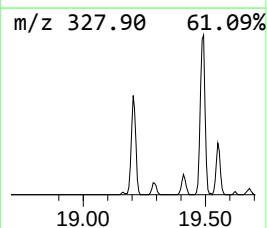
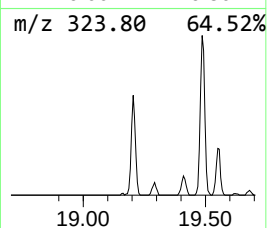
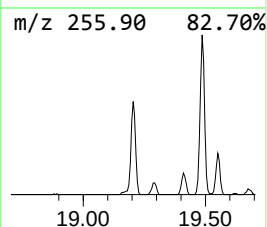
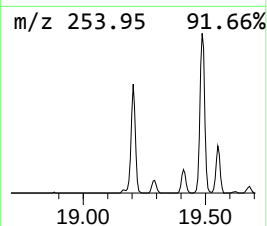
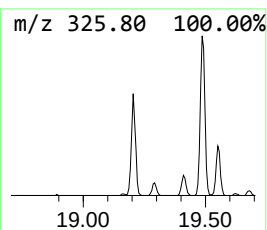
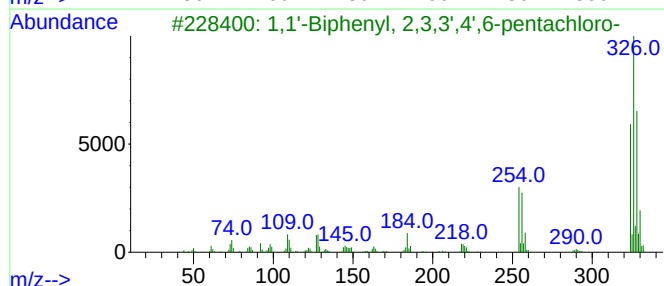
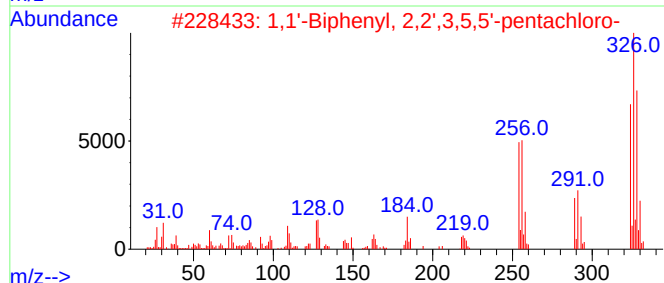
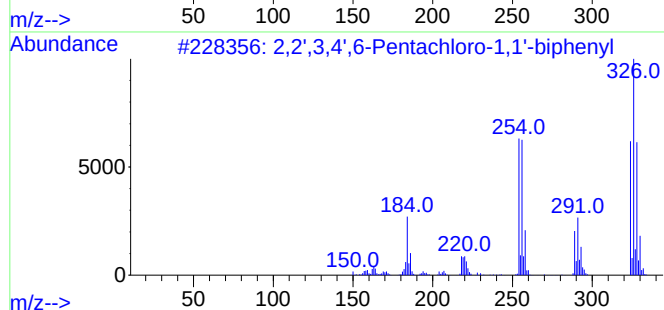
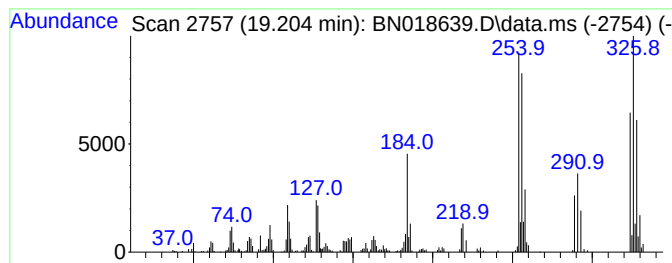
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.204	3.80 ng/ul	387578	Phenanthrene-d10	17.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4',6-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-05-8	99		
2	1,1'-Biphenyl, 2,2',3,5,5'-penta...	324	C12H5Cl5	052663-61-3	99		
3	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99		
4	2,3,3',4',5'- Pentachloro-1,1'-b...	324	C12H5Cl5	076842-07-4	99		
5	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	99		



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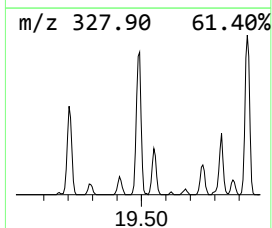
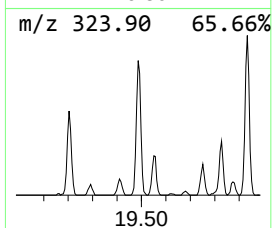
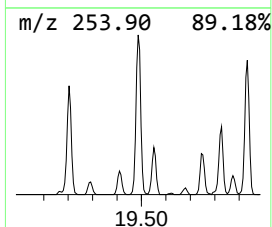
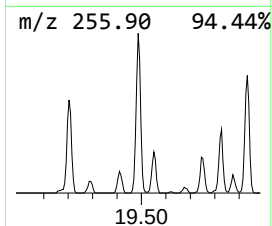
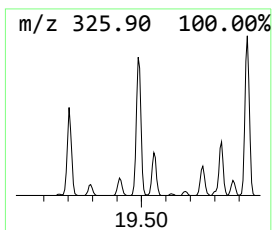
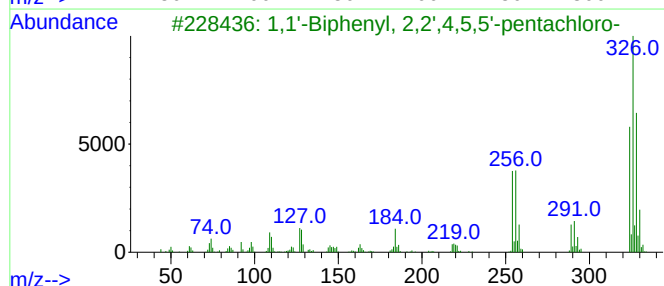
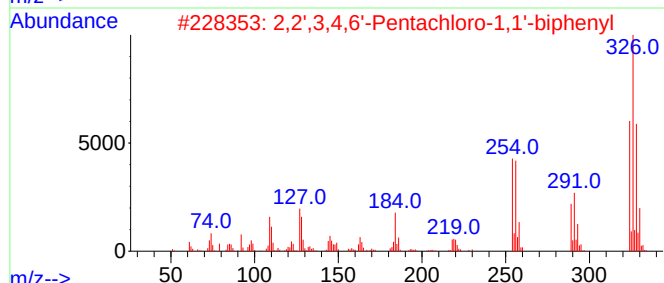
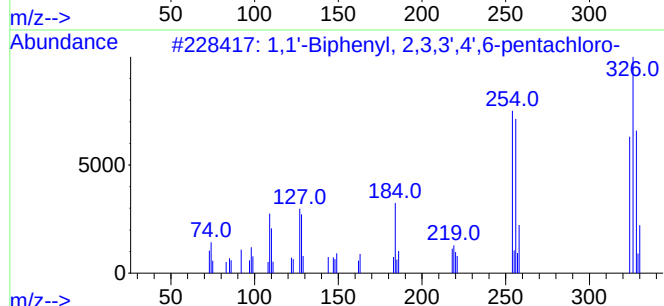
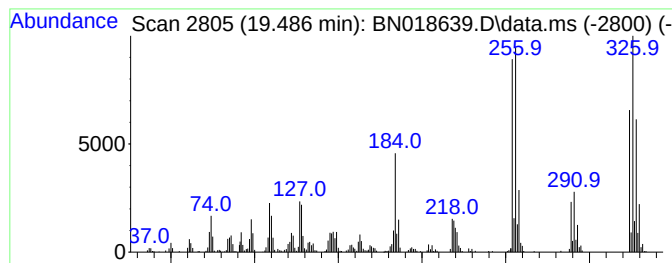
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Peak Number 4 1,1'-Biphenyl, 2,3,3',4',6-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.486	6.45 ng/ul	581038	Chrysene-d12	21.463		
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	2,2',3,4,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-57-2	99	
3	1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	99	
4	3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	99	
5	2,2',3,6,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-54-9	99	



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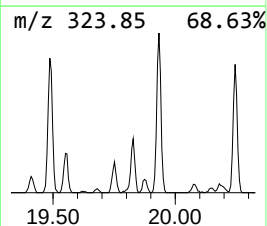
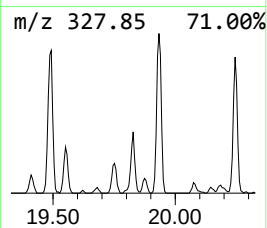
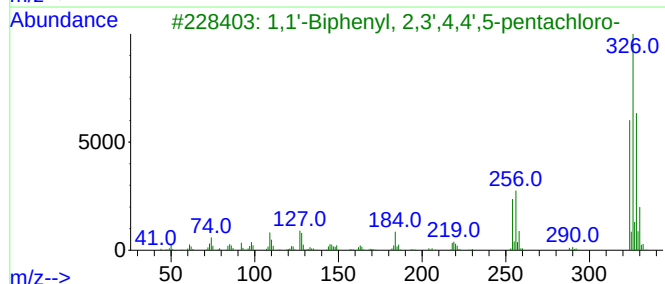
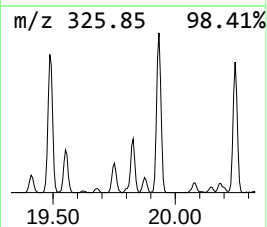
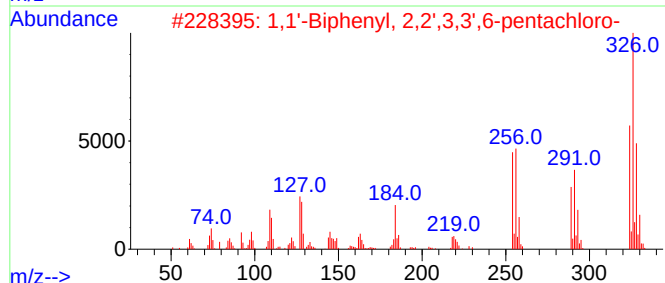
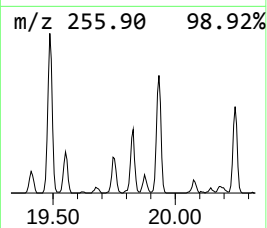
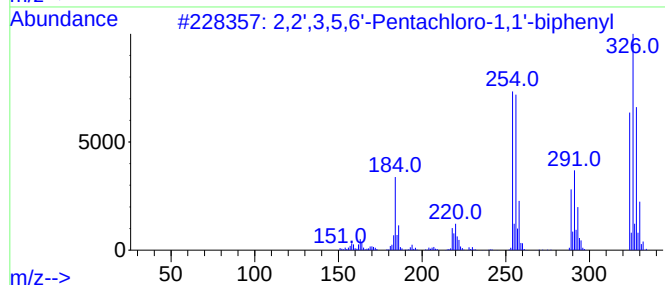
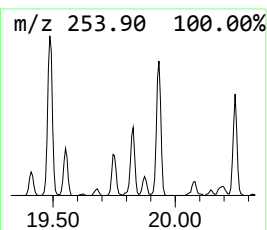
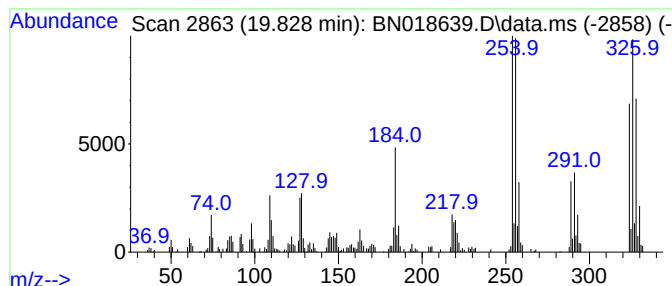
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.828	2.43 ng/ul	219140	Chrysene-d12	21.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,5,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-55-0	99		
2	1,1'-Biphenyl, 2,2',3,3',6-penta...	324	C12H5Cl5	052663-60-2	99		
3	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99		
4	3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	99		
5	1,1'-Biphenyl, 2,2',3',4,5-Penta...	324	C12H5Cl5	041464-51-1	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021722\
Data File : BN018639.D
Acq On : 17 Feb 2022 22:04
Operator : CG/JU
Sample : N1507-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

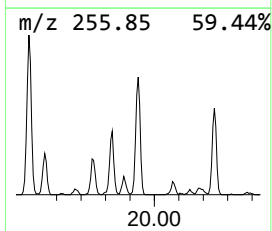
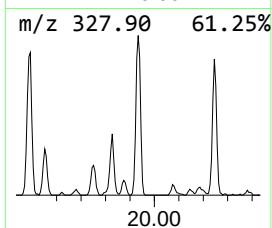
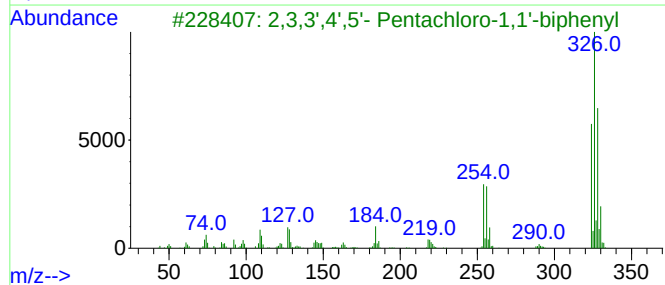
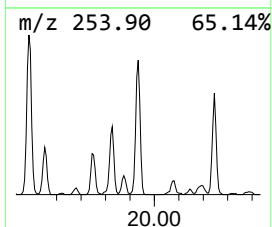
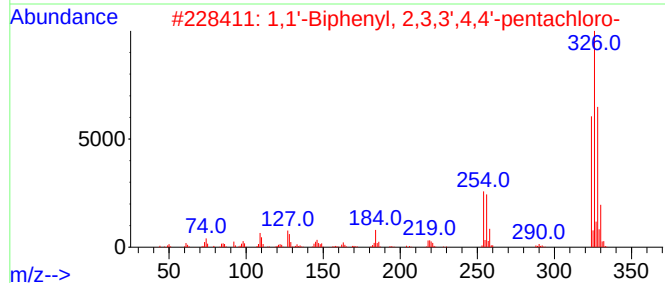
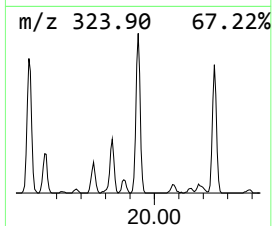
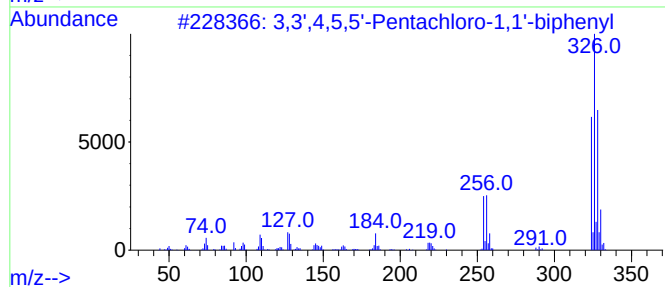
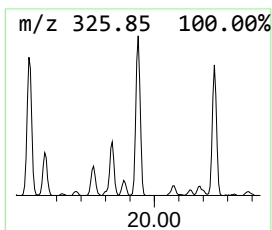
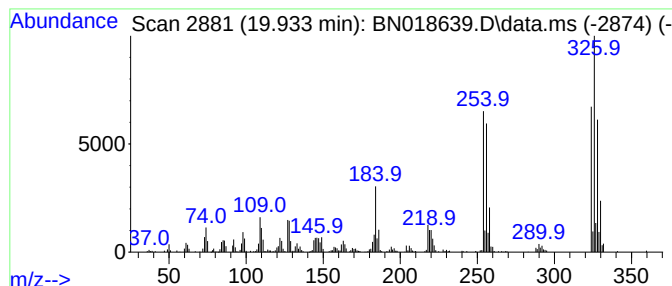
Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 3,3',4,5,5'-Pentachloro-1,1... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.933	5.73 ng/ul	516621	Chrysene-d12	21.463	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3',4,5,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-33-1	99
2	1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99
3	2,3,3',4',5'- Pentachloro-1,1'-b...	324	C12H5Cl5	076842-07-4	99
4	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	99
5	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021722\
Data File : BN018639.D
Acq On : 17 Feb 2022 22:04
Operator : CG/JU
Sample : N1507-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

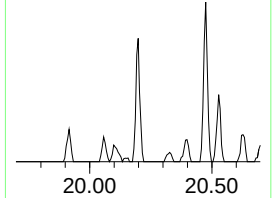
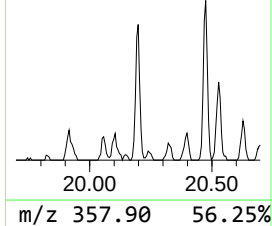
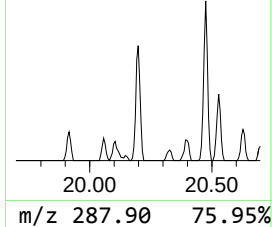
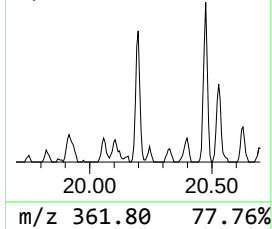
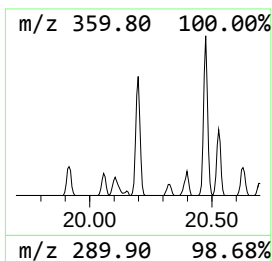
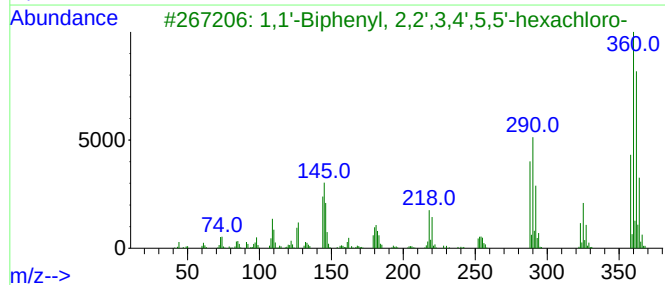
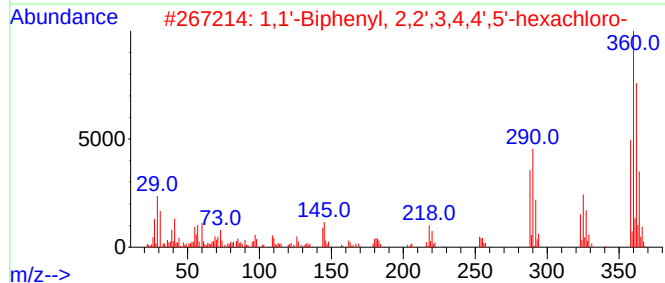
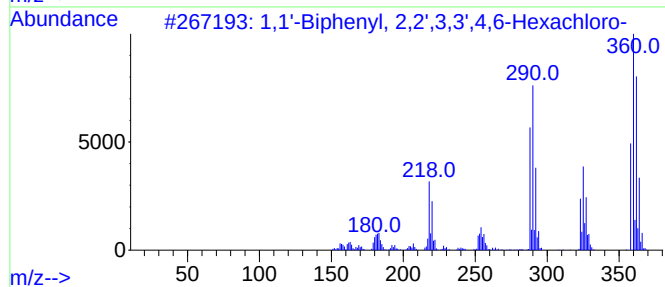
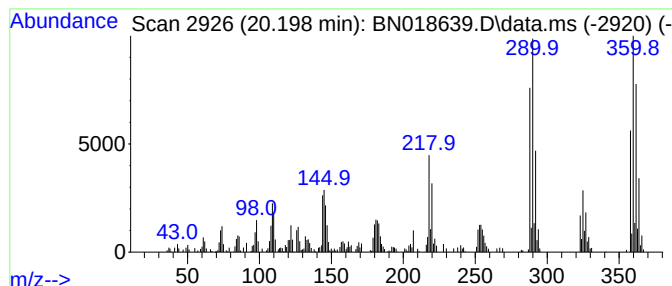
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.198	2.78 ng/ul	250079	Chrysene-d12	21.463	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,3',4,6-Hex...	358	C12H4Cl6	061798-70-7	99
2	1,1'-Biphenyl, 2,2',3,4,4',5'-he...	358	C12H4Cl6	035065-28-2	99
3	1,1'-Biphenyl, 2,2',3,4',5,5'-he...	358	C12H4Cl6	051908-16-8	99
4	1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99
5	1,1'-Biphenyl, 2,2',4,4',5',6-He...	358	C12H4Cl6	060145-22-4	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021722\
Data File : BN018639.D
Acq On : 17 Feb 2022 22:04
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Sample : N1507-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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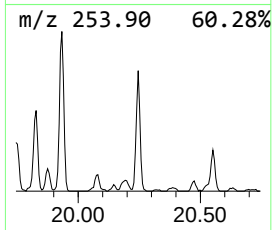
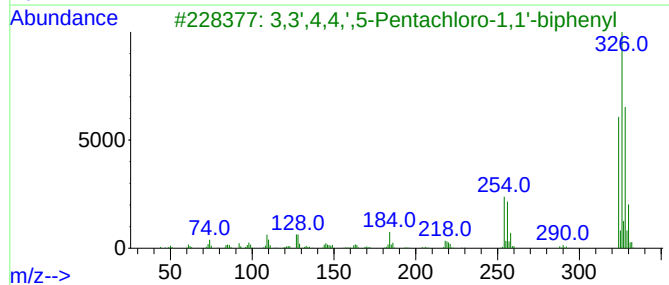
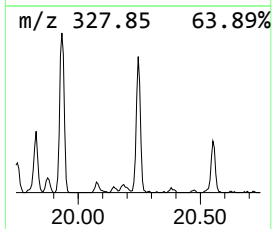
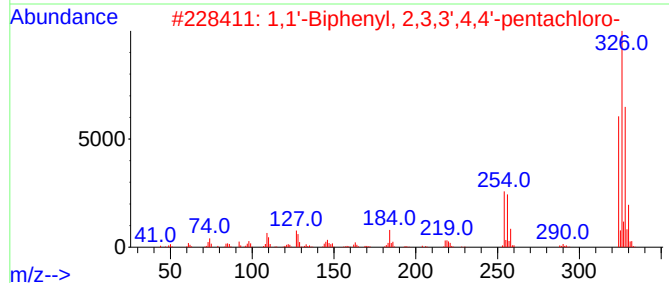
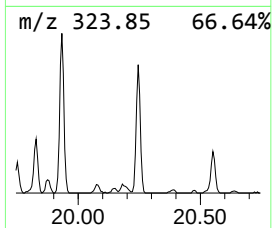
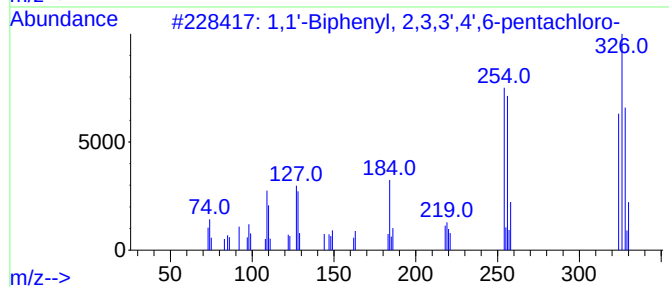
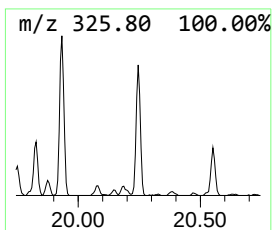
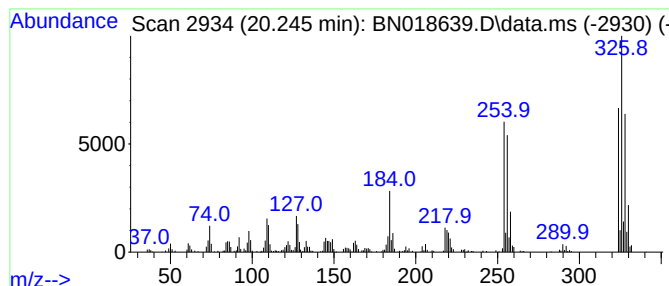
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1,1'-Biphenyl, 2,3,3',4,4'-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.245	3.92 ng/ul	352832	Chrysene-d12	21.463

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, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99		
3	3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	99		
4	1,1'-Biphenyl, 2,2',3',4,5-Penta...	324	C12H5Cl5	041464-51-1	99		
5	3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021722\
Data File : BN018639.D
Acq On : 17 Feb 2022 22:04
Operator : CG/JU
Sample : N1507-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
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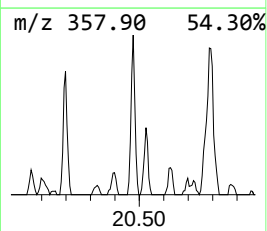
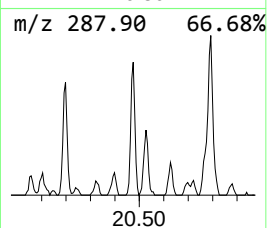
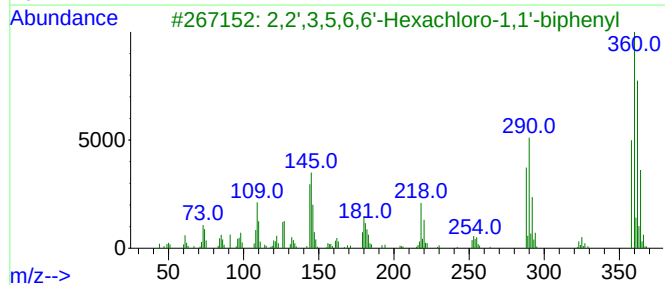
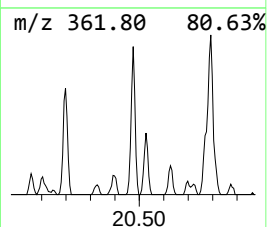
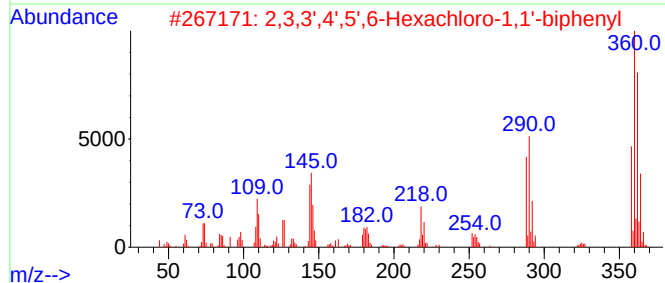
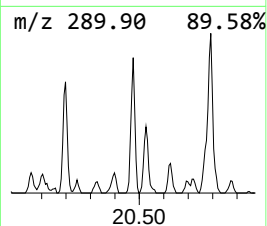
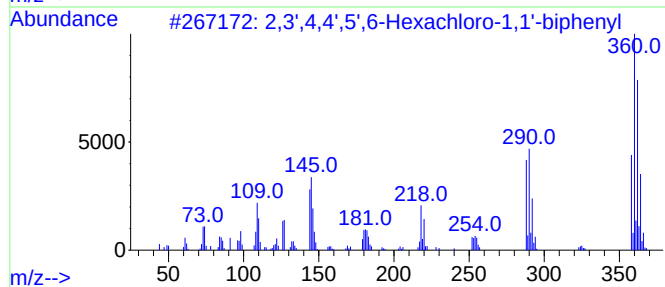
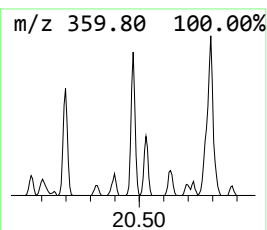
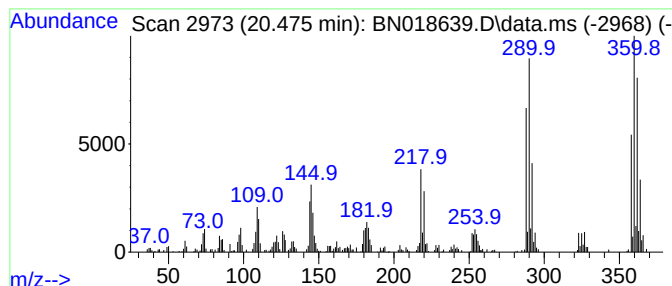
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2,3',4,4',5',6-Hexachloro-1... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.474	2.72 ng/ul	245478	Chrysene-d12	21.463

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',4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-45-0	99		
3	2,2',3,5,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-09-2	99		
4	2,2',3,4',6,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	068194-08-1	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\BN021722\
Data File : BN018639.D
Acq On : 17 Feb 2022 22:04
Operator : CG/JU
Sample : N1507-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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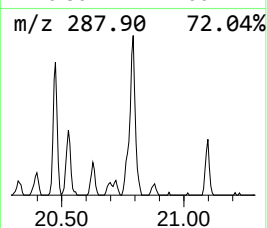
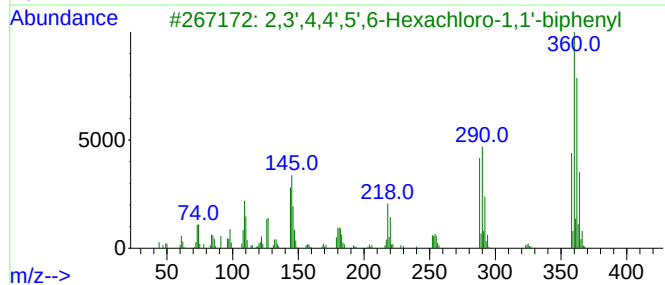
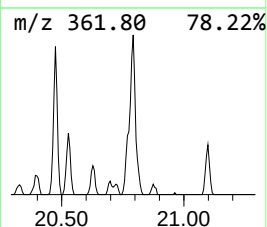
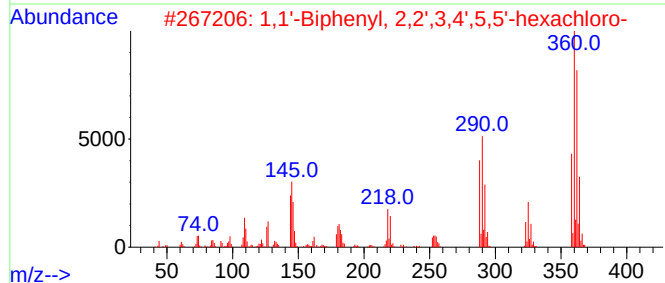
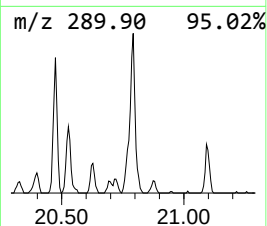
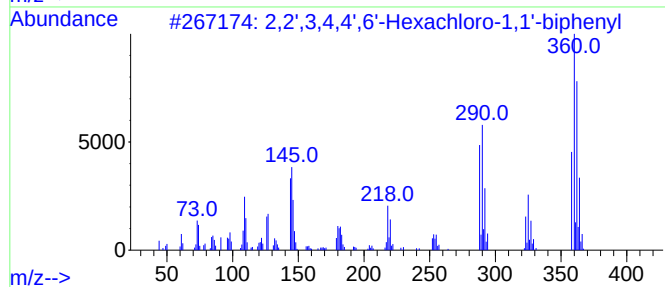
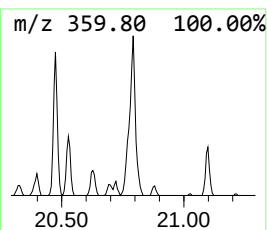
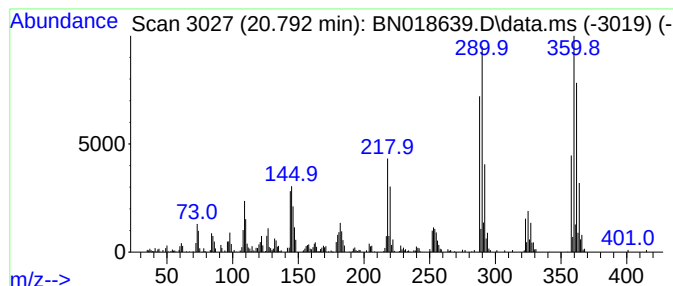
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.792	4.47 ng/ul	402325	Chrysene-d12	21.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4,4',6'-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-64-4	99		
2	1,1'-Biphenyl, 2,2',3,4',5,5'-he...	358	C12H4Cl6	051908-16-8	99		
3	2,3',4,4',5',6'-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	99		
4	1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	99		
5	1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021722\
Data File : BN018639.D
Acq On : 17 Feb 2022 22:04
Operator : CG/JU
Sample : N1507-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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(DEL) Alkane: S...	3.223	56.1	ng/ul	3652990	1	7.916	1303330	20.0
2,2',3,4',6-Pen...	19.204	3.8	ng/ul	387578	4	17.287	2039610	20.0
1,1'-Biphenyl, ...	19.486	6.5	ng/ul	581038	5	21.463	1802030	20.0
2,2',3,5,6'-Pen...	19.828	2.4	ng/ul	219140	5	21.463	1802030	20.0
3,3',4,5,5'-Pen...	19.933	5.7	ng/ul	516621	5	21.463	1802030	20.0
1,1'-Biphenyl, ...	20.198	2.8	ng/ul	250079	5	21.463	1802030	20.0
1,1'-Biphenyl, ...	20.245	3.9	ng/ul	352832	5	21.463	1802030	20.0
2,3',4,4',5',6-...	20.474	2.7	ng/ul	245478	5	21.463	1802030	20.0
2,2',3,4,4',6'-...	20.792	4.5	ng/ul	402325	5	21.463	1802030	20.0