

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021722\
 Data File : BN018637.D
 Acq On : 17 Feb 2022 20:52
 Operator : CG/JU
 Sample : N1507-10
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGZF3

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: BN018637.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	16	rVB	216874	241446	6.46%	1.201%
2	3.123	17	23	26	rBV2	63709	74288	1.99%	0.369%
3	3.152	26	28	30	rVV	26429	24221	0.65%	0.120%
4	3.181	30	33	35	rVV	38340	36162	0.97%	0.180%
5	3.223	35	40	44	rVV	3594904	3739100	100.00%	18.596%
6	3.540	90	94	95	rBV2	19237	19766	0.53%	0.098%
7	3.570	95	99	104	rVB	139841	166303	4.45%	0.827%
8	3.664	112	115	120	rVB5	9707	12804	0.34%	0.064%
9	3.964	162	166	174	rVV2	36309	53206	1.42%	0.265%
10	4.064	177	183	188	rVB3	42329	68601	1.83%	0.341%
11	4.622	272	278	283	rVB2	17204	26182	0.70%	0.130%
12	4.928	326	330	336	rVB2	10119	14123	0.38%	0.070%
13	4.999	336	342	346	rBV6	7588	11975	0.32%	0.060%
14	5.358	396	403	405	rBV5	5311	10395	0.28%	0.052%
15	5.405	407	411	416	rVB	12848	18011	0.48%	0.090%
16	5.540	429	434	439	rBV3	7149	11312	0.30%	0.056%
17	5.822	478	482	487	rBV2	14070	19636	0.53%	0.098%
18	5.911	494	497	503	rVB7	4809	8989	0.24%	0.045%
19	6.187	536	544	550	rVB5	6660	11840	0.32%	0.059%
20	6.399	575	580	585	rVB2	13564	21644	0.58%	0.108%
21	7.305	729	734	738	rBV4	9168	14942	0.40%	0.074%
22	7.346	738	741	745	rVB2	8766	12006	0.32%	0.060%
23	7.399	745	750	751	rBV3	7830	10440	0.28%	0.052%
24	7.487	760	765	770	rBV3	9471	15707	0.42%	0.078%
25	7.546	770	775	776	rBV3	8258	8860	0.24%	0.044%
26	7.916	831	838	849	rBV	816577	1308209	34.99%	6.506%
27	8.058	855	862	867	rVB4	10313	18085	0.48%	0.090%
28	8.210	883	888	896	rVB7	5255	11276	0.30%	0.056%
29	8.469	927	932	937	rVB6	6190	11932	0.32%	0.059%
30	8.593	950	953	960	rVB3	6396	10244	0.27%	0.051%
31	8.699	964	971	973	rBV6	6031	9728	0.26%	0.048%
32	9.163	1040	1050	1061	rVB	28053	56208	1.50%	0.280%
33	9.293	1068	1072	1077	rVB7	7252	13451	0.36%	0.067%
34	9.910	1172	1177	1185	rVB9	5994	11492	0.31%	0.057%
35	10.163	1216	1220	1229	rVB7	6317	13899	0.37%	0.069%
36	10.722	1304	1315	1321	rBV	1052462	1749380	46.79%	8.700%

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
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37	10.769	1321	1323	1329	rVV	35331	52854	1.41%	0.263%
38	10.869	1333	1340	1342	rBV6	5562	10316	0.28%	0.051%
39	10.899	1342	1345	1351	rVB3	6163	9423	0.25%	0.047%
40	10.993	1355	1361	1367	rVB8	10850	20018	0.54%	0.100%
41	11.204	1391	1397	1404	rVB3	13508	24960	0.67%	0.124%
42	11.428	1430	1435	1440	rVV5	9500	16005	0.43%	0.080%
43	11.504	1442	1448	1452	rBV8	7643	13475	0.36%	0.067%
44	12.157	1554	1559	1565	rVB2	9589	15641	0.42%	0.078%
45	12.381	1590	1597	1603	rVB2	10174	18281	0.49%	0.091%
46	12.598	1629	1634	1643	rVB	7533	13062	0.35%	0.065%
47	13.387	1763	1768	1778	rVB2	20805	32871	0.88%	0.163%
48	14.045	1875	1880	1883	rVB5	6586	9270	0.25%	0.046%
49	14.387	1934	1938	1945	rVB6	4444	9408	0.25%	0.047%
50	14.457	1945	1950	1956	rBV2	14493	22505	0.60%	0.112%
51	14.551	1956	1966	1974	rBV	1338164	2011067	53.78%	10.002%
52	14.945	2027	2033	2038	rVB	14550	20178	0.54%	0.100%
53	15.404	2106	2111	2115	rVB	12102	14836	0.40%	0.074%
54	15.592	2138	2143	2148	rBV3	15186	22230	0.59%	0.111%
55	16.022	2211	2216	2222	rBV8	10034	21661	0.58%	0.108%
56	16.263	2253	2257	2259	rBV2	18296	22839	0.61%	0.114%
57	16.287	2259	2261	2265	rVB4	13119	16987	0.45%	0.084%
58	16.516	2296	2300	2305	rBV8	5168	10831	0.29%	0.054%
59	16.598	2308	2314	2318	rBV	23038	37071	0.99%	0.184%
60	17.051	2388	2391	2393	rBV2	10665	13236	0.35%	0.066%
61	17.104	2398	2400	2407	rVB5	16782	23976	0.64%	0.119%
62	17.169	2408	2411	2415	rBV5	8974	14523	0.39%	0.072%
63	17.286	2424	2431	2436	rBV	1442217	2085814	55.78%	10.374%
64	17.328	2436	2438	2443	rVV	182031	213616	5.71%	1.062%
65	17.392	2447	2449	2454	rVB5	16668	21596	0.58%	0.107%
66	17.498	2459	2467	2468	rBV7	13223	33789	0.90%	0.168%
67	17.645	2489	2492	2496	rBV5	33701	49466	1.32%	0.246%
68	17.857	2525	2528	2532	rVB4	44072	50260	1.34%	0.250%
69	18.039	2555	2559	2562	rBV2	57004	76966	2.06%	0.383%
70	18.210	2585	2588	2592	rVB2	42636	48103	1.29%	0.239%
71	18.263	2592	2597	2606	rVB5	104807	183727	4.91%	0.914%
72	18.357	2609	2613	2618	rVB2	162428	214236	5.73%	1.065%
73	18.416	2621	2623	2628	rVB3	26533	32016	0.86%	0.159%
74	18.463	2628	2631	2634	rBV4	21196	35544	0.95%	0.177%
75	18.645	2658	2662	2670	rVB3	65196	117547	3.14%	0.585%
76	18.769	2679	2683	2688	rBV2	55267	82874	2.22%	0.412%

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77	18.910	2703	2707	2712	rBV	131485	154166	4.12%	0.767%
78	19.033	2724	2728	2732	rVB2	39515	59972	1.60%	0.298%
79	19.180	2749	2753	2754	rBV	76259	97155	2.60%	0.483%
80	19.204	2754	2757	2762	rVB2	207243	280092	7.49%	1.393%
81	19.269	2765	2768	2770	rBV2	27772	38913	1.04%	0.194%
82	19.339	2776	2780	2786	rBV	186438	243108	6.50%	1.209%
83	19.404	2787	2791	2796	rVV2	84204	129772	3.47%	0.645%
84	19.486	2800	2805	2812	rVV	285046	373069	9.98%	1.855%
85	19.551	2812	2816	2820	rVV2	78377	104892	2.81%	0.522%
86	19.592	2820	2823	2831	rVB	46549	76932	2.06%	0.383%
87	19.745	2846	2849	2854	rBV3	62441	78920	2.11%	0.393%
88	19.828	2859	2863	2866	rBV	99031	120734	3.23%	0.600%
89	19.869	2867	2870	2875	rVV5	56331	83884	2.24%	0.417%
90	19.933	2877	2881	2885	rVB	248036	301955	8.08%	1.502%
91	20.198	2922	2926	2929	rBV2	127197	161238	4.31%	0.802%
92	20.245	2930	2934	2938	rVB	167270	203915	5.45%	1.014%
93	20.475	2970	2973	2978	rVB3	120583	161443	4.32%	0.803%
94	20.527	2979	2982	2984	rBV3	62778	75257	2.01%	0.374%
95	20.627	2996	2999	3003	rVB	100913	107226	2.87%	0.533%
96	20.722	3012	3015	3020	rVB2	61482	80794	2.16%	0.402%
97	20.786	3024	3026	3033	rVB	126641	162785	4.35%	0.810%
98	21.463	3136	3141	3145	rBV	1324734	1697698	45.40%	8.443%
99	21.745	3186	3189	3192	rBV2	85958	94256	2.52%	0.469%
100	23.863	3542	3549	3557	rVB2	777445	1609590	43.05%	8.005%

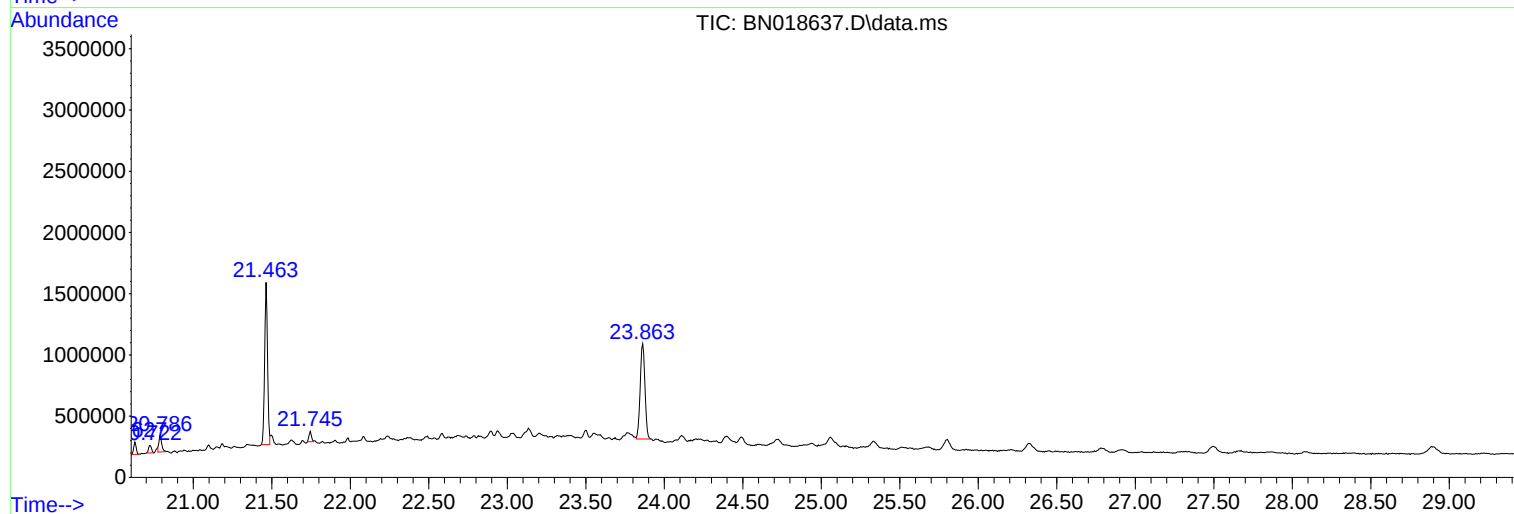
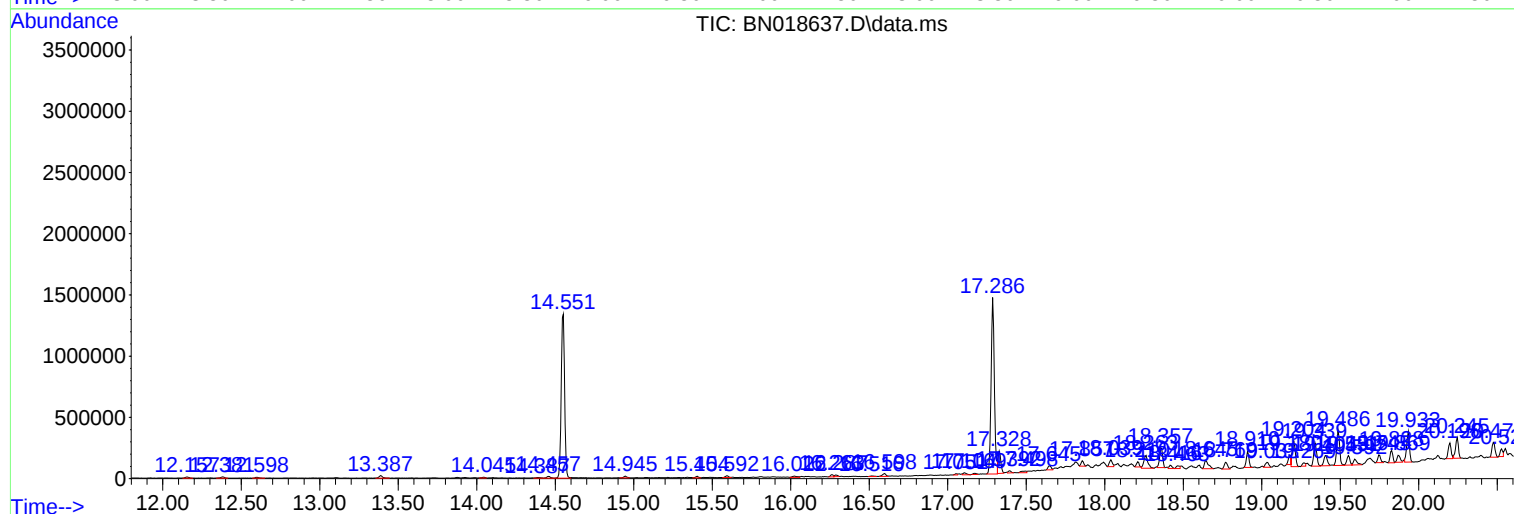
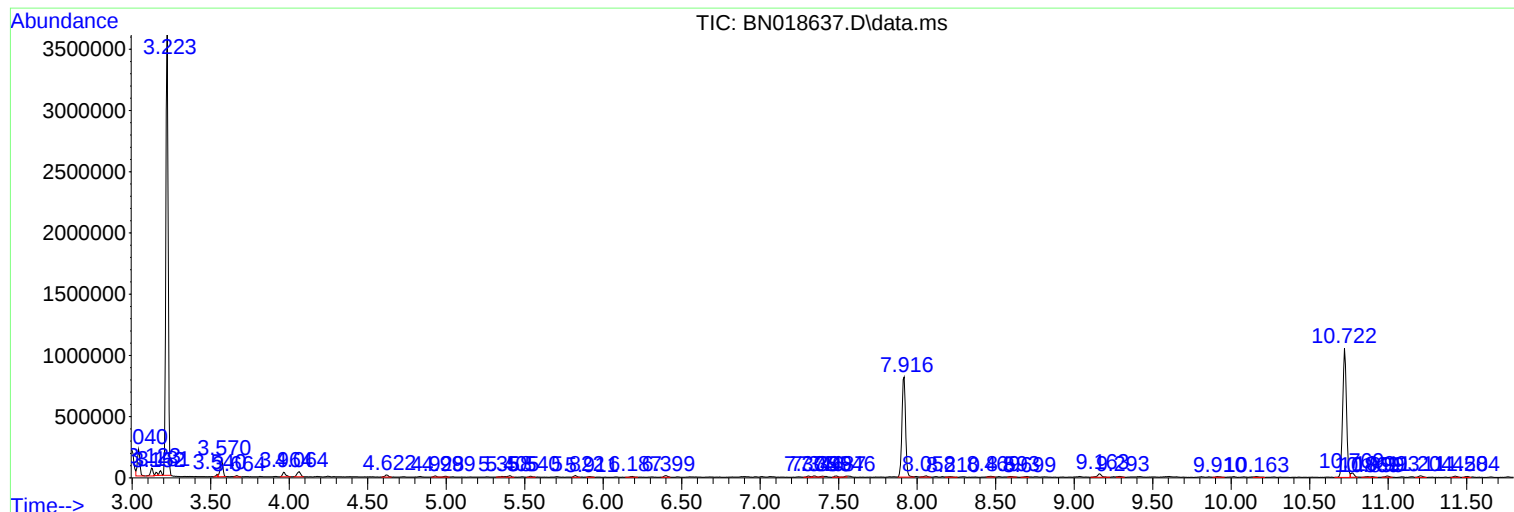
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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



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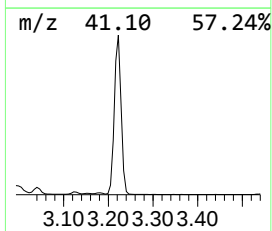
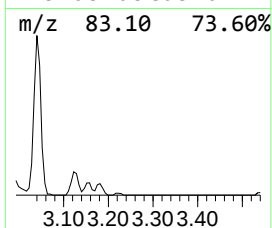
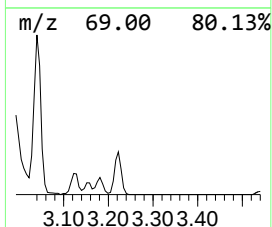
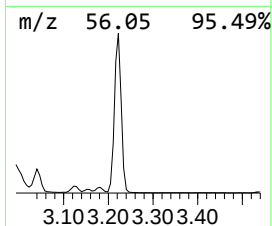
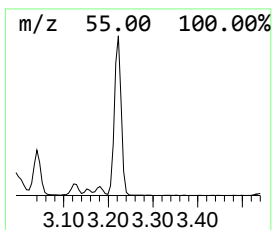
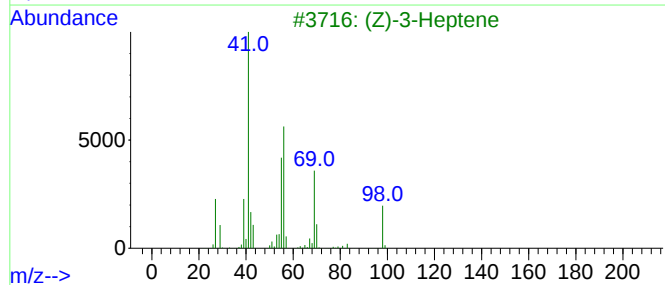
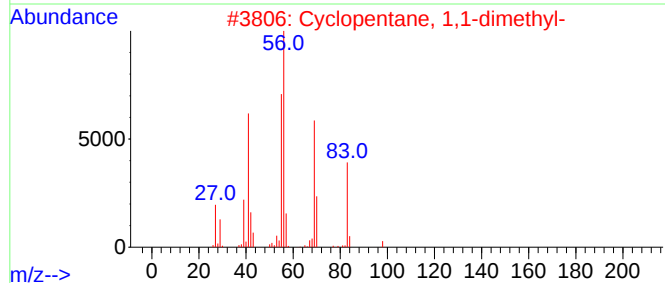
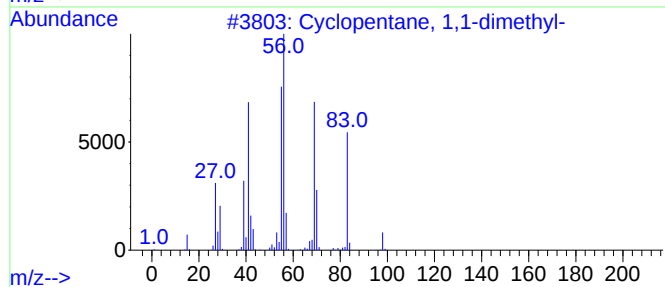
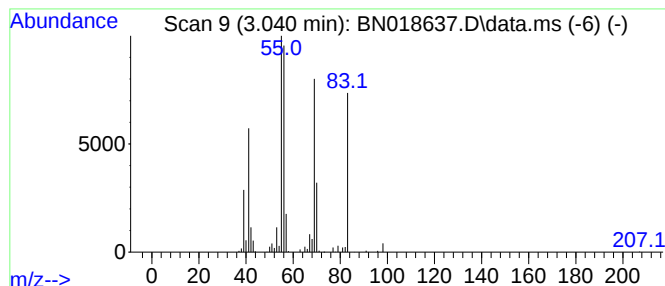
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.040 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.040	3.69 ng/ul	241446	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			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	58
3			(Z)-3-Heptene	98	C7H14	007642-10-6	53
4			3-Heptene	98	C7H14	000592-78-9	53
5			3-Heptene, (E)-	98	C7H14	014686-14-7	50



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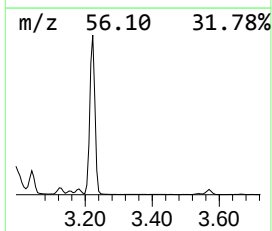
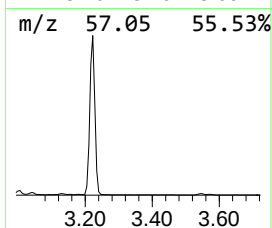
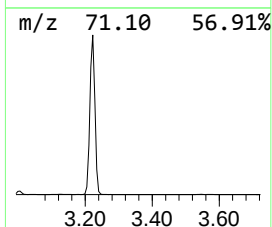
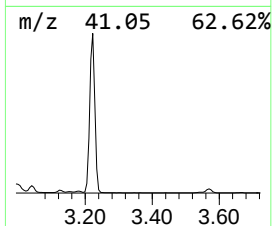
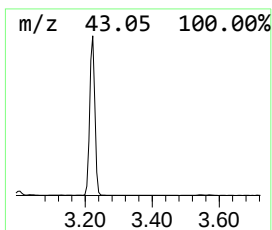
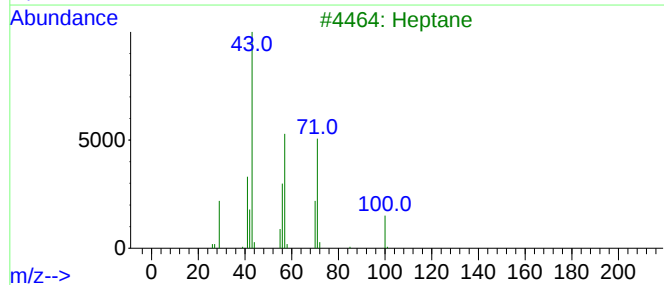
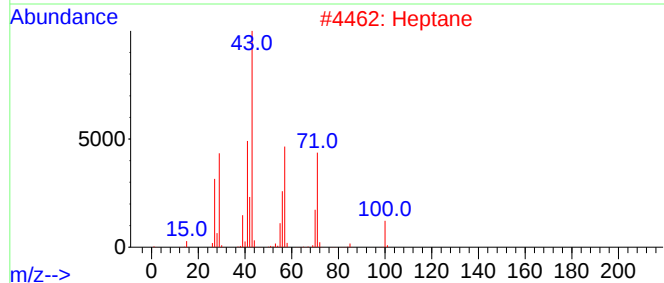
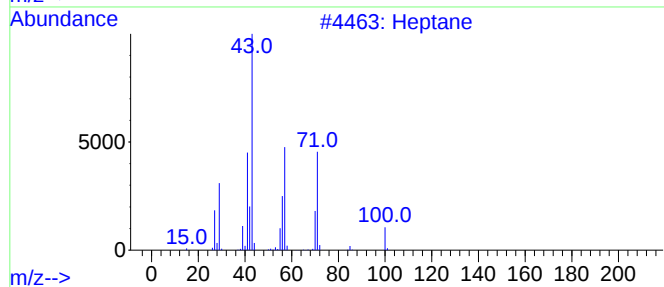
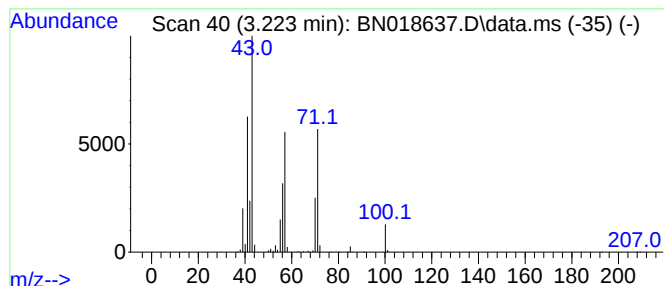
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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	57.16 ng/ul	3739100	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	94
3	Heptane			100	C7H16	000142-82-5	91
4	Heptane			100	C7H16	000142-82-5	87
5	Hexane, 3-methyl-			100	C7H16	000589-34-4	45



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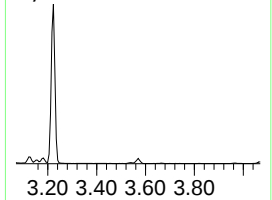
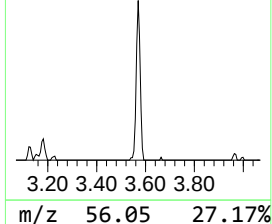
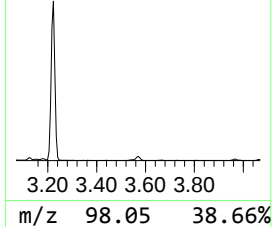
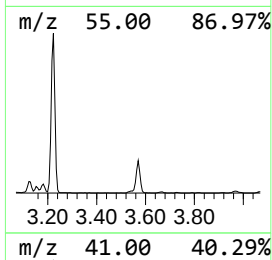
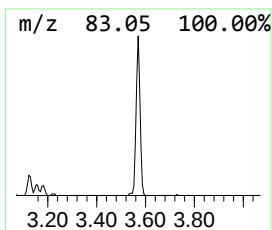
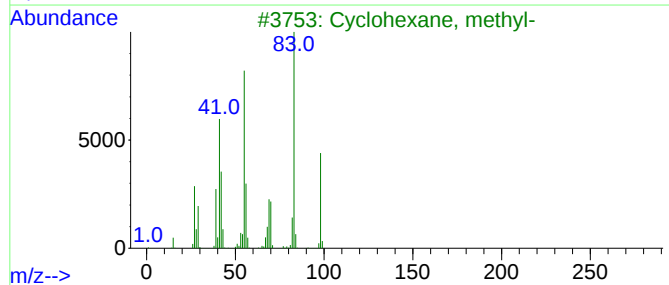
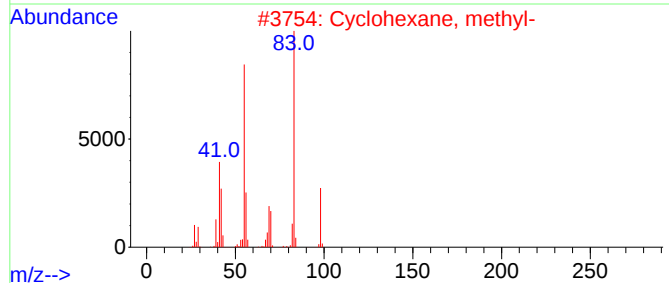
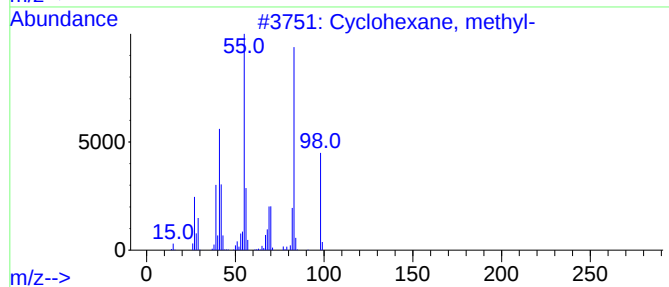
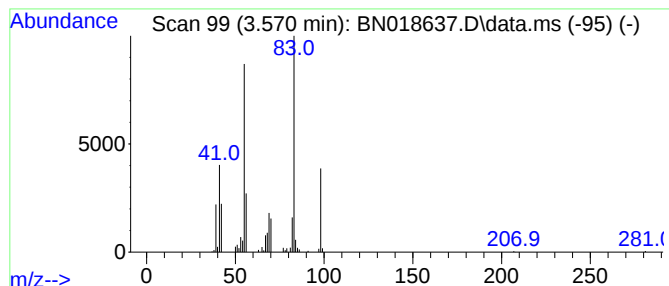
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Cyclic3.570 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.570	2.54 ng/ul	166303	1,4-Dichlorobenzene-d4	7.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, methyl-	98	C7H14	000108-87-2	94
2			Cyclohexane, methyl-	98	C7H14	000108-87-2	91
3			Cyclohexane, methyl-	98	C7H14	000108-87-2	87
4			Cyclohexane, methyl-	98	C7H14	000108-87-2	87
5			Cyclohexane, methyl-	98	C7H14	000108-87-2	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021722\
Data File : BN018637.D
Acq On : 17 Feb 2022 20:52
Operator : CG/JU
Sample : N1507-10
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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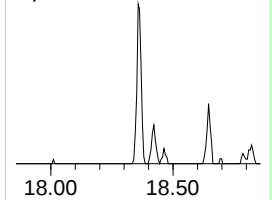
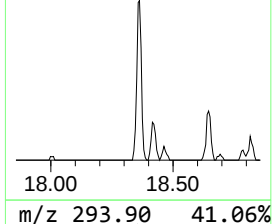
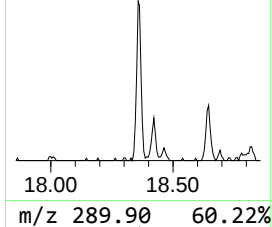
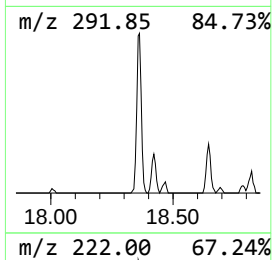
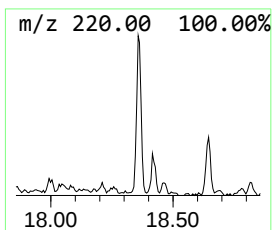
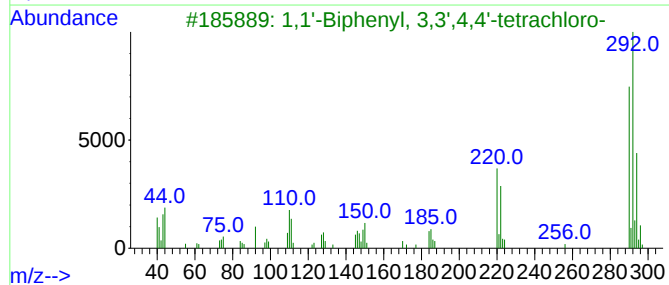
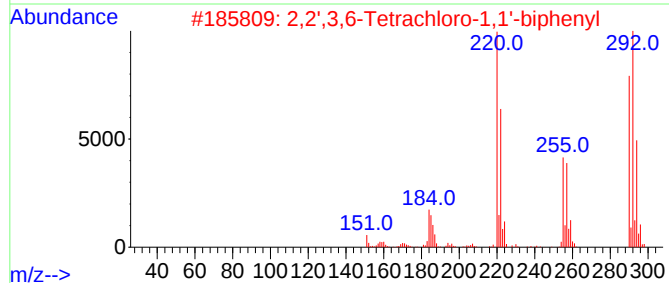
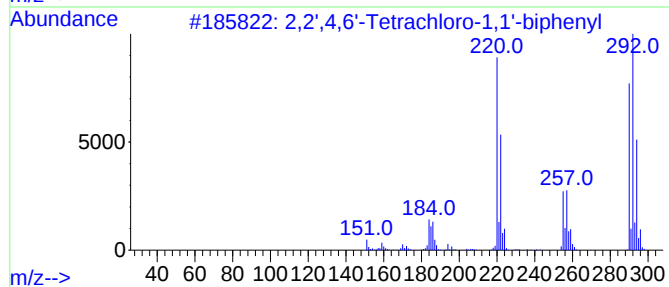
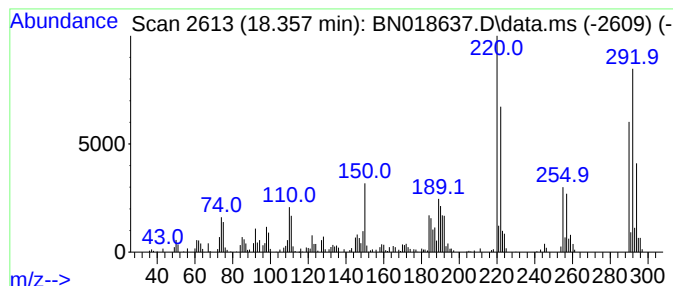
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2,2',4,6'-Tetrachloro-1,1'-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.357	2.05 ng/ul	214236	Phenanthrene-d10	17.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99		
2	2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	99		
3	1,1'-Biphenyl, 3,3',4,4'-tetrachloro-	290	C12H6Cl4	032598-13-3	99		
4	1,1'-Biphenyl, 3,3',4,5'-tetrachloro-	290	C12H6Cl4	041464-48-6	99		
5	1,1'-Biphenyl, 2,2',4,4'-tetrachloro-	290	C12H6Cl4	002437-79-8	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021722\
Data File : BN018637.D
Acq On : 17 Feb 2022 20:52
Operator : CG/JU
Sample : N1507-10
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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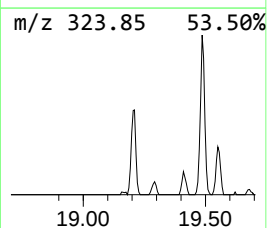
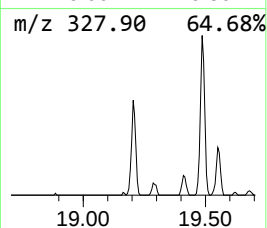
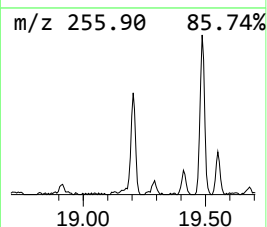
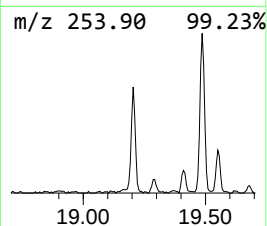
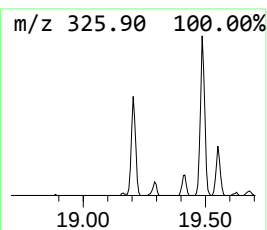
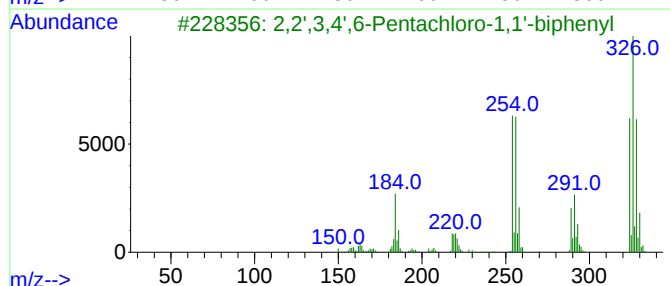
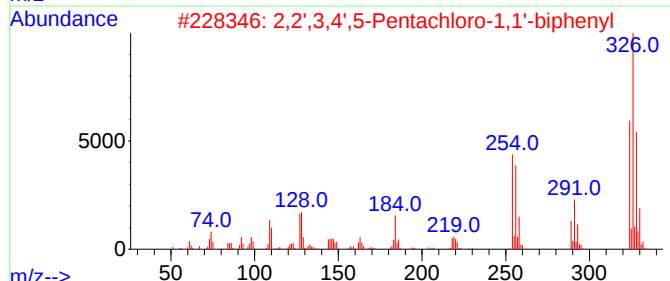
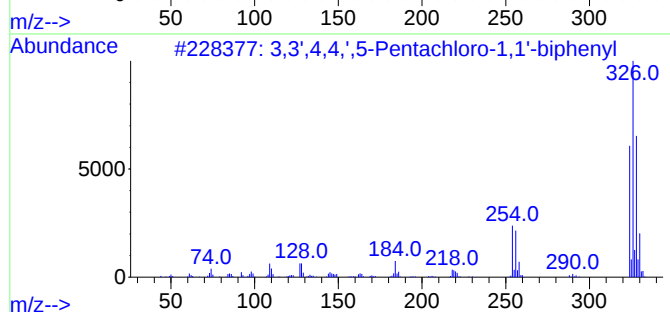
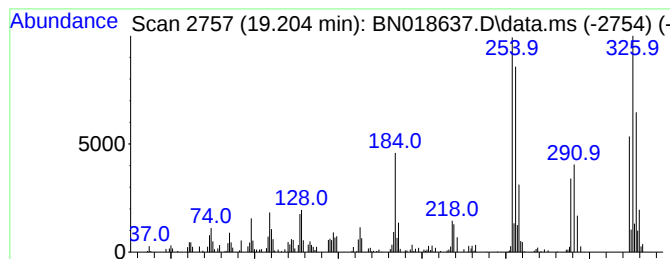
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.204	2.69 ng/ul	280092	Phenanthrene-d10	17.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	99		
2	2,2',3,4',5-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-07-0	99		
3	2,2',3,4',6-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-05-8	99		
4	1,1'-Biphenyl, 2,2',3,4,5'-penta...	324	C12H5Cl5	038380-02-8	99		
5	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021722\
Data File : BN018637.D
Acq On : 17 Feb 2022 20:52
Operator : CG/JU
Sample : N1507-10
Misc :
ALS Vial : 12 Sample Multiplier: 1

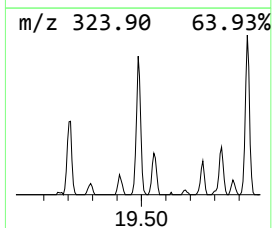
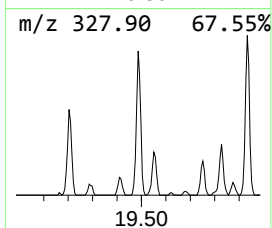
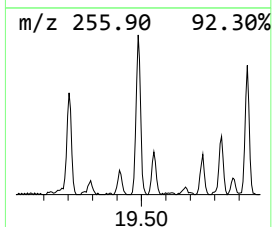
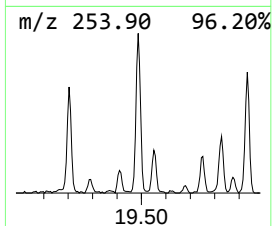
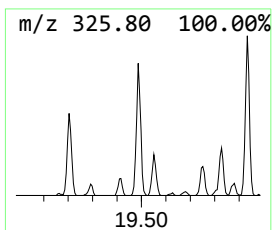
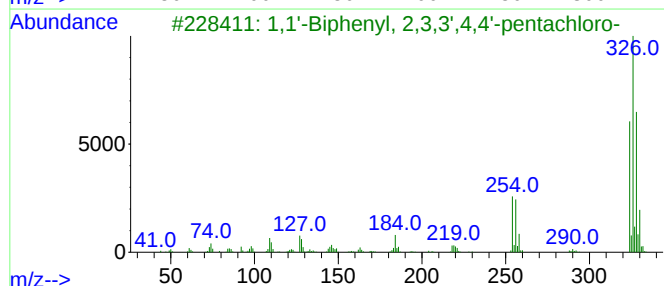
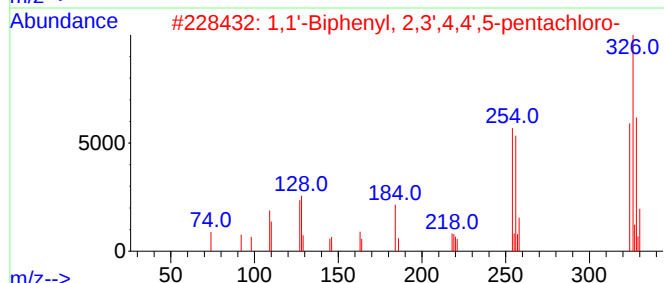
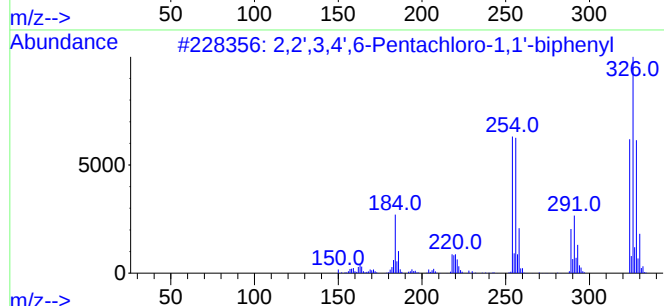
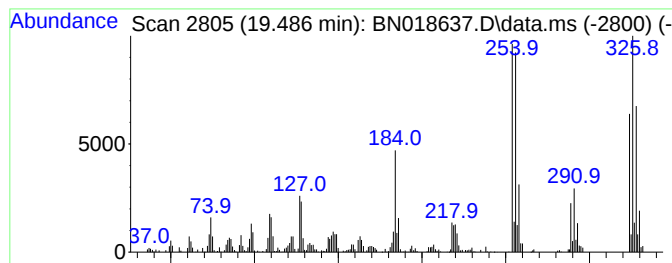
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.486	4.39 ng/ul	373069	Chrysene-d12	21.463	
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, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99
3	1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99
4	2,2',3,6,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-54-9	99
5	2,3,3',4,5-Pentachloro-1,1'-biph...	324	C12H5Cl5	070424-69-0	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021722\
Data File : BN018637.D
Acq On : 17 Feb 2022 20:52
Operator : CG/JU
Sample : N1507-10
Misc :
ALS Vial : 12 Sample Multiplier: 1

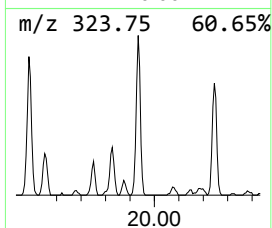
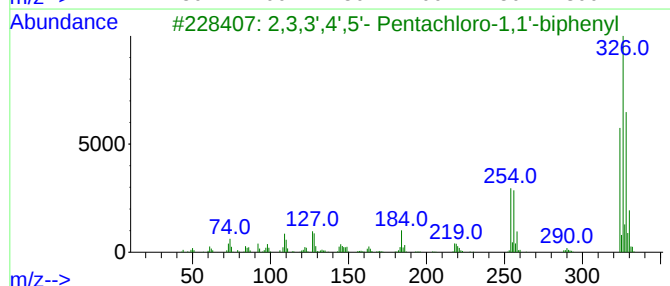
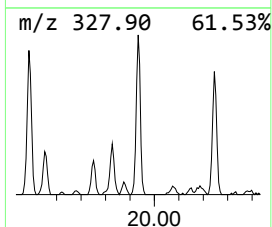
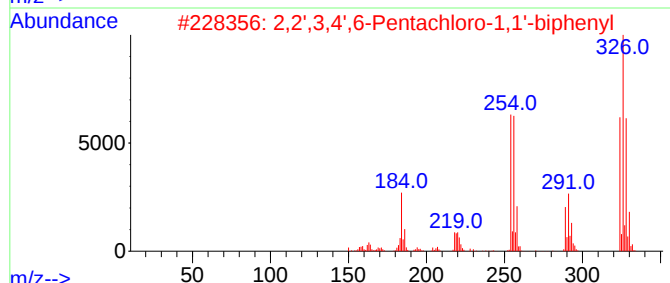
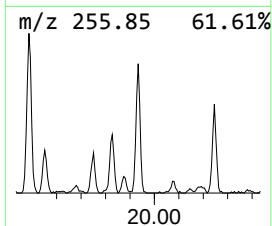
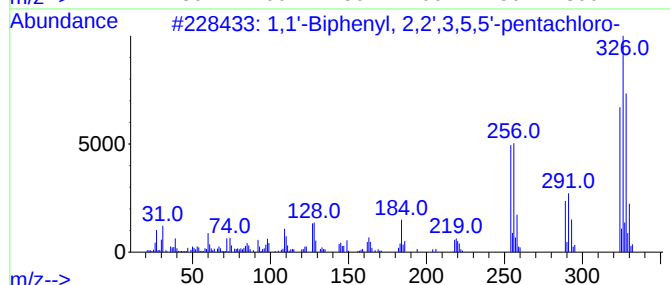
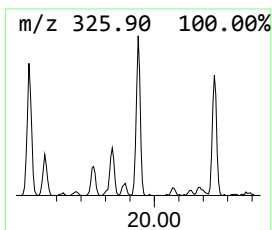
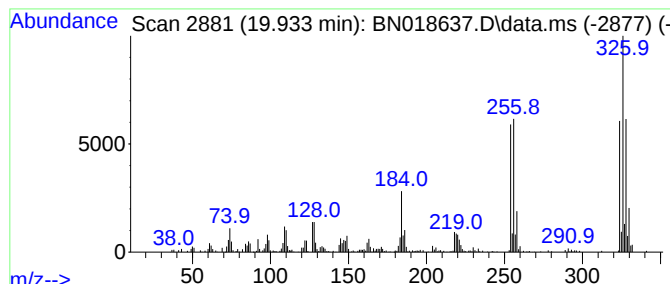
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.933	3.56 ng/ul	301955	Chrysene-d12	21.463		
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,5,5'-penta...	324	C12H5Cl5	052663-61-3	99	
2	2,2',3,4',6-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-05-8	99	
3	2,3,3',4',5'- Pentachloro-1,1'-b...	324	C12H5Cl5	076842-07-4	99	
4	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	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\
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Acq On : 17 Feb 2022 20:52
Operator : CG/JU
Sample : N1507-10
Misc :
ALS Vial : 12 Sample Multiplier: 1

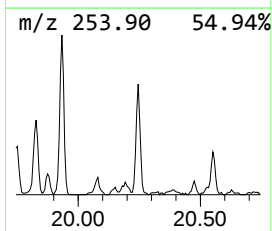
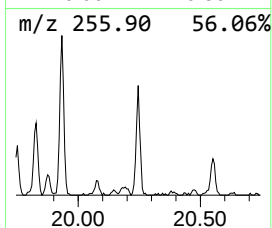
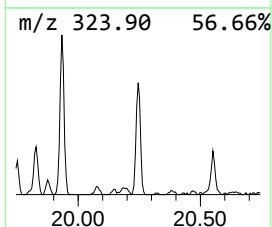
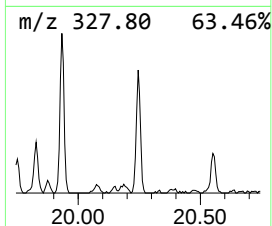
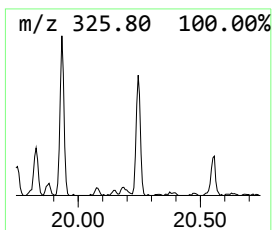
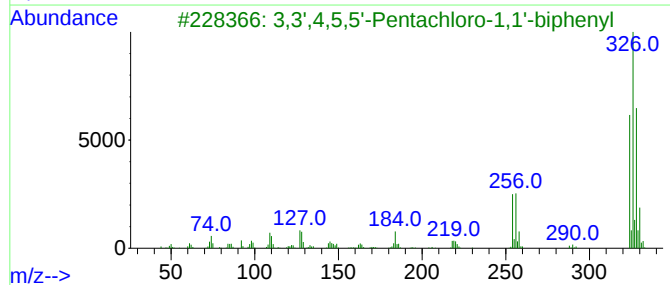
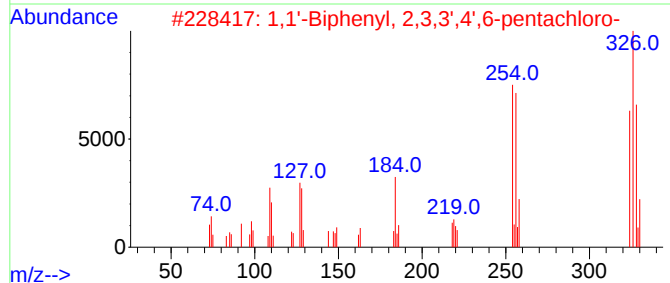
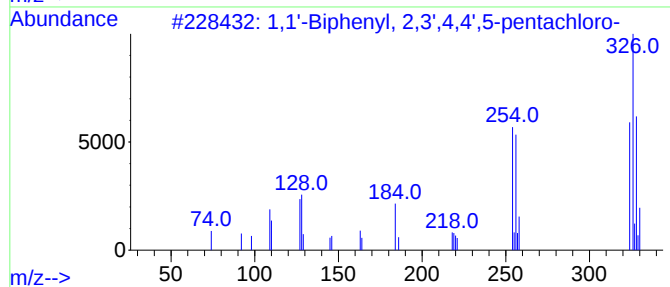
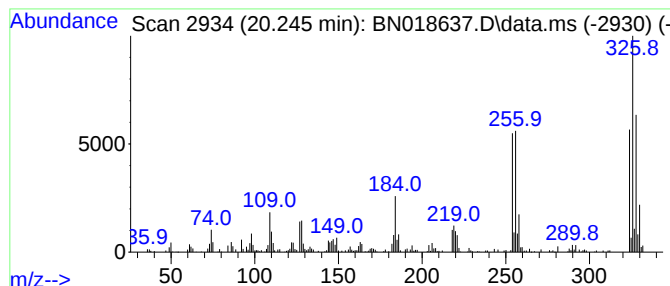
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 1,1'-Biphenyl, 2,3',4,4',5-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.245	2.40 ng/ul	203915	Chrysene-d12	21.463	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99
2	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99
3	3,3',4,5,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-33-1	99
4	2,3,3',4,5-Pentachloro-1,1'-biph...	324	C12H5Cl5	070424-69-0	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 : BN018637.D
Acq On : 17 Feb 2022 20:52
Operator : CG/JU
Sample : N1507-10
Misc :
ALS Vial : 12 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\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	57.2	ng/ul	3739100	1	7.916	1308210	20.0
(DEL) Alkane: C...	3.570	2.5	ng/ul	166303	1	7.916	1308210	20.0
2,2',4,6'-Tetra...	18.357	2.0	ng/ul	214236	4	17.287	2085810	20.0
3,3',4,4',5-Pe...	19.204	2.7	ng/ul	280092	4	17.287	2085810	20.0
2,2',3,4',6-Pen...	19.486	4.4	ng/ul	373069	5	21.463	1697700	20.0
1,1'-Biphenyl, ...	19.933	3.6	ng/ul	301955	5	21.463	1697700	20.0
1,1'-Biphenyl, ...	20.245	2.4	ng/ul	203915	5	21.463	1697700	20.0