

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
 Data File : BP010163.D
 Acq On : 29 Apr 2022 22:19
 Operator : CG/JU
 Sample : N2543-01
 Misc : GCMS CONFIRMATION
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETNS5

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_P\Methods\SFAM-EPA-BP042122.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010163.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.117	3	5	11	rVB	192986	205991	5.93%	1.014%
2	3.199	16	19	22	rVV	60050	60497	1.74%	0.298%
3	3.228	22	24	26	rVV	25470	25319	0.73%	0.125%
4	3.252	26	28	31	rVV	36196	40470	1.17%	0.199%
5	3.293	31	35	40	rVB	3408744	3473285	100.00%	17.092%
6	3.599	84	87	88	rBV3	6273	5916	0.17%	0.029%
7	3.628	88	92	99	rVB	86857	110579	3.18%	0.544%
8	3.723	104	108	115	rVB8	6688	10305	0.30%	0.051%
9	4.105	168	173	179	rVB	17625	25857	0.74%	0.127%
10	4.558	248	250	255	rVB2	4652	5411	0.16%	0.027%
11	5.234	362	365	370	rVB7	2742	4156	0.12%	0.020%
12	7.922	815	822	835	rBV	747173	1213210	34.93%	5.970%
13	10.510	1257	1262	1264	rBV6	2760	4439	0.13%	0.022%
14	10.728	1288	1299	1312	rBV	892963	1611641	46.40%	7.931%
15	10.869	1318	1323	1330	rVB9	5662	12740	0.37%	0.063%
16	11.210	1379	1381	1387	rVB7	3342	4636	0.13%	0.023%
17	13.398	1745	1753	1760	rVB7	3027	6115	0.18%	0.030%
18	14.557	1943	1950	1962	rBV	1357017	2001605	57.63%	9.850%
19	14.728	1975	1979	1983	rVB6	5577	8024	0.23%	0.039%
20	15.269	2066	2071	2076	rBV9	1894	4273	0.12%	0.021%
21	15.351	2080	2085	2089	rBV7	1922	3808	0.11%	0.019%
22	15.581	2118	2124	2126	rBV6	2350	4861	0.14%	0.024%
23	15.816	2160	2164	2176	rVB6	12291	22977	0.66%	0.113%
24	16.839	2333	2338	2343	rVB2	46339	63835	1.84%	0.314%
25	17.192	2392	2398	2401	rBV	257704	360998	10.39%	1.776%
26	17.228	2401	2404	2411	rVB	58203	77037	2.22%	0.379%
27	17.310	2411	2418	2433	rBV	1442315	2295694	66.10%	11.297%
28	17.486	2443	2448	2460	rVB2	165666	259309	7.47%	1.276%
29	17.763	2491	2495	2499	rBV	37851	50655	1.46%	0.249%
30	17.916	2513	2521	2527	rBV	326211	565393	16.28%	2.782%
31	17.981	2530	2532	2535	rVV3	11428	13099	0.38%	0.064%
32	18.028	2535	2540	2547	rVB5	37324	64863	1.87%	0.319%
33	18.098	2549	2552	2556	rVV6	10657	12418	0.36%	0.061%
34	18.151	2556	2561	2566	rVV	120215	167456	4.82%	0.824%
35	18.204	2566	2570	2575	rVB4	40303	56195	1.62%	0.277%
36	18.310	2585	2588	2592	rVB6	17816	21827	0.63%	0.107%

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Integrator: RTE

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37	18.363	2592	2597	2602	rVV	111229	157147	4.52%	0.773%
38	18.422	2603	2607	2611	rVV	87052	113996	3.28%	0.561%
39	18.463	2611	2614	2620	rVB	75242	97434	2.81%	0.479%
40	18.639	2639	2644	2648	rBV	124909	174277	5.02%	0.858%
41	18.686	2649	2652	2656	rVV2	49687	62253	1.79%	0.306%
42	18.733	2656	2660	2664	rVV	68392	99638	2.87%	0.490%
43	18.775	2664	2667	2670	rVV4	45515	63653	1.83%	0.313%
44	18.810	2670	2673	2680	rVB	90003	114043	3.28%	0.561%
45	18.904	2686	2689	2697	rVB4	32454	43534	1.25%	0.214%
46	19.104	2719	2723	2727	rBV	57098	74743	2.15%	0.368%
47	19.157	2727	2732	2735	rVV	110355	146718	4.22%	0.722%
48	19.192	2735	2738	2747	rVB	113805	148619	4.28%	0.731%
49	19.398	2768	2773	2779	rBV3	84538	174622	5.03%	0.859%
50	19.451	2779	2782	2787	rVB7	20043	27202	0.78%	0.134%
51	19.775	2833	2837	2842	rVB	231047	268198	7.72%	1.320%
52	19.875	2850	2854	2860	rVB	166740	200640	5.78%	0.987%
53	20.269	2916	2921	2925	rBV	280140	320183	9.22%	1.576%
54	20.322	2926	2930	2936	rVB	166848	202472	5.83%	0.996%
55	20.698	2989	2994	2999	rBV	1240665	1401912	40.36%	6.899%
56	21.392	3107	3112	3125	rBV	1356627	1931322	55.61%	9.504%
57	23.839	3522	3528	3543	rVB	712709	1624149	46.76%	7.992%

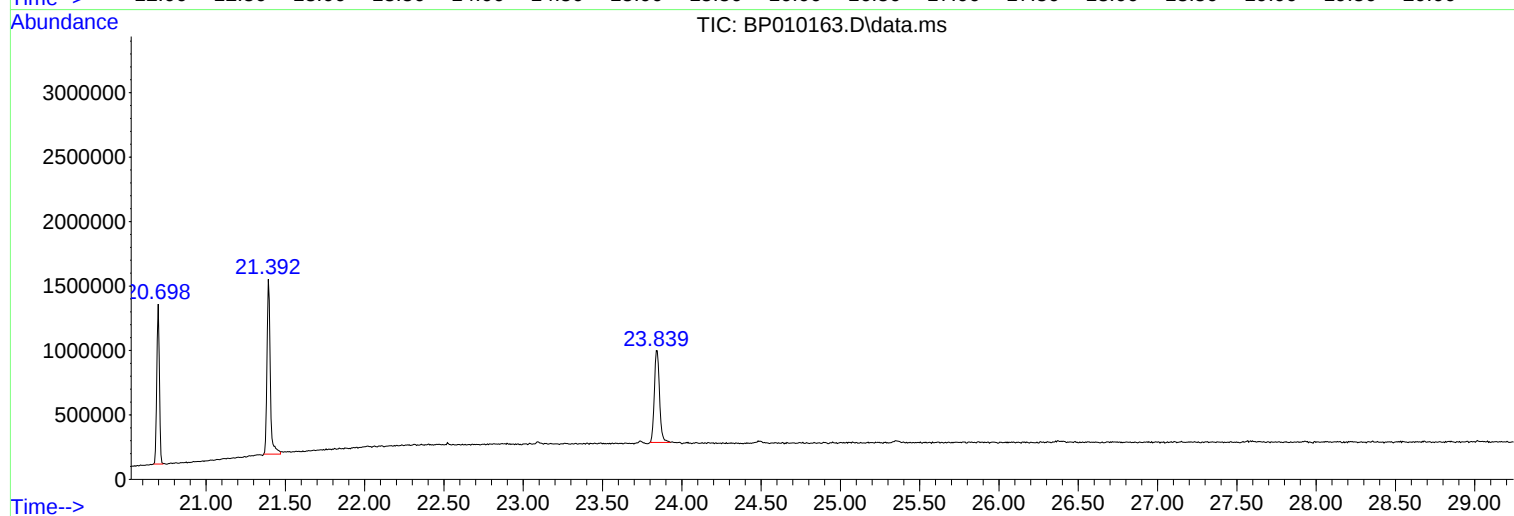
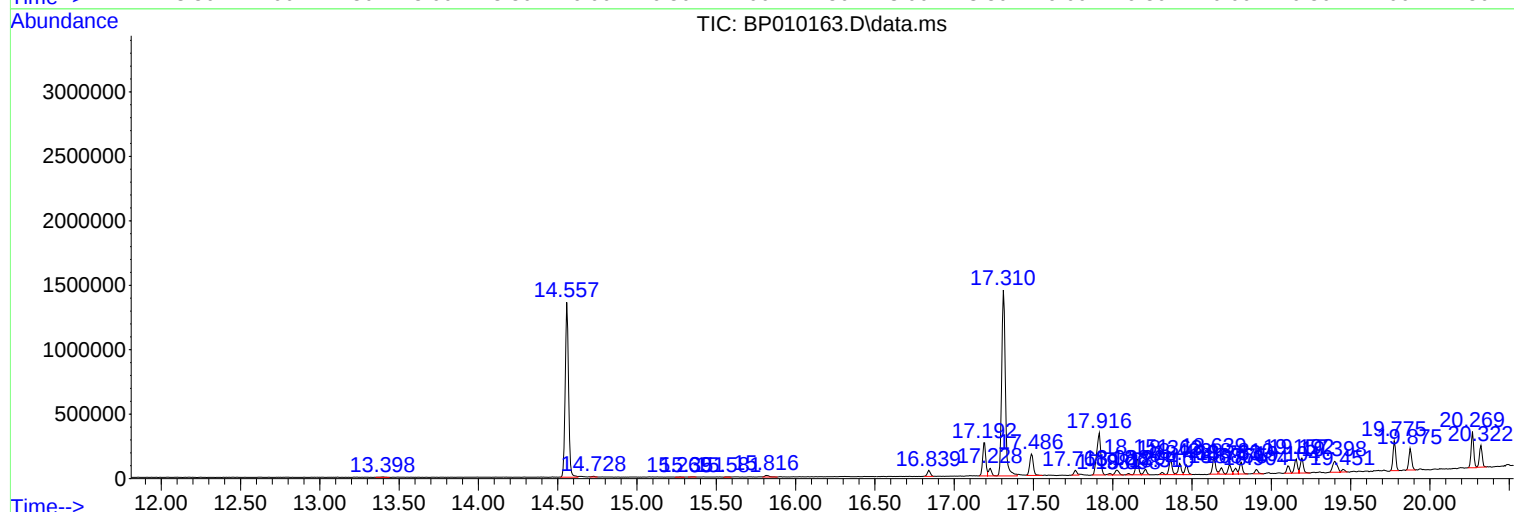
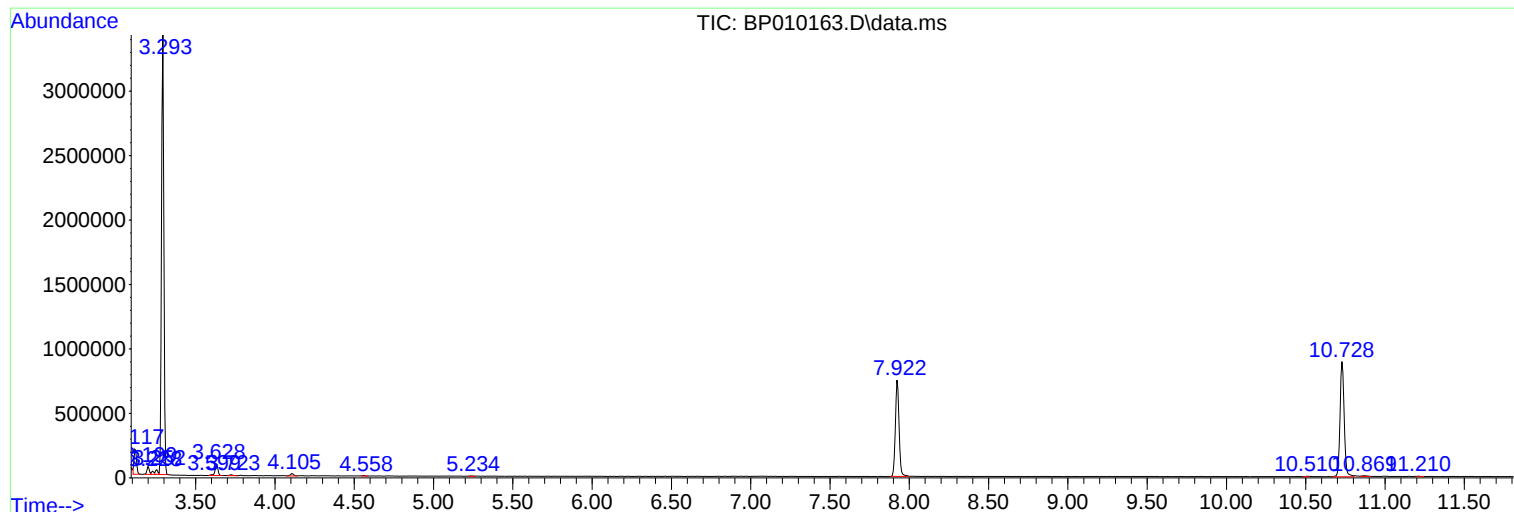
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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



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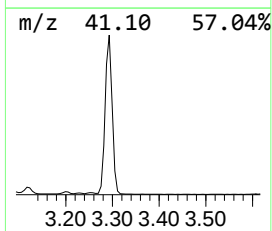
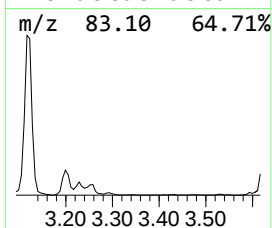
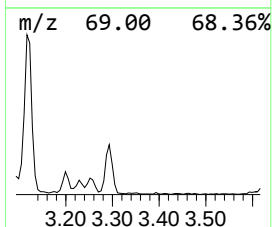
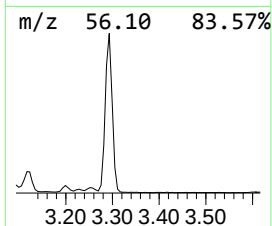
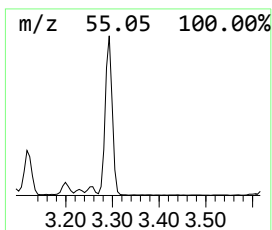
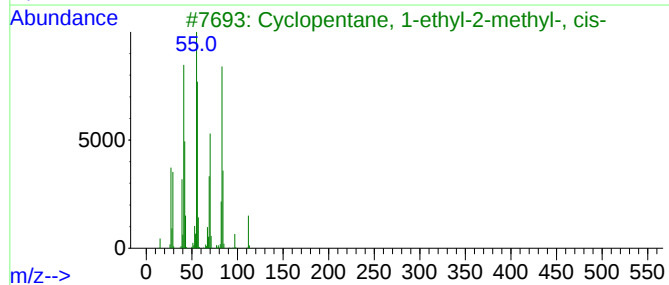
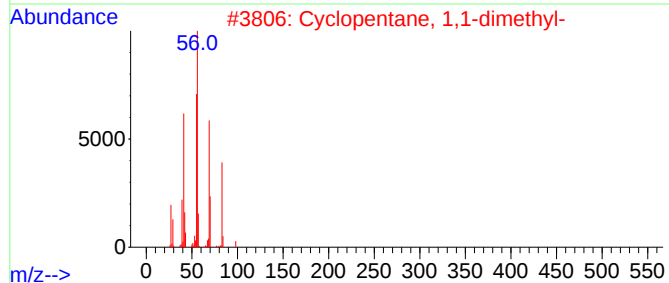
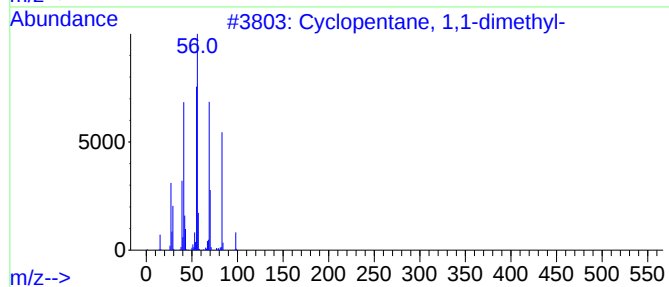
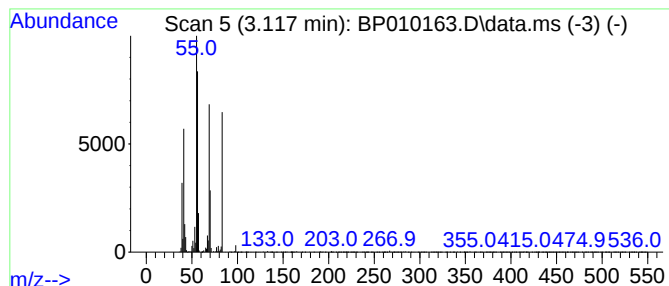
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.117 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.117	3.40 ng/ul	205991	1,4-Dichlorobenzene-d4	7.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	78
2			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	64
3			Cyclopentane, 1-ethyl-2-methyl-,...	112	C8H16	000930-89-2	59
4			3-Heptene	98	C7H14	000592-78-9	50
5			2-Hexene, 5-methyl-, (E)-	98	C7H14	007385-82-2	50



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ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
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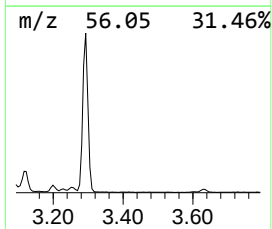
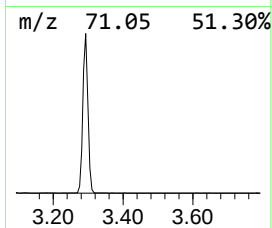
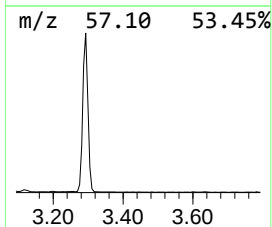
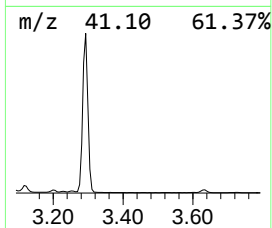
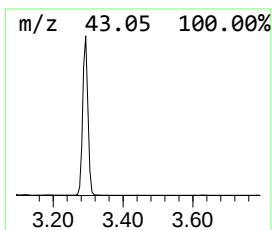
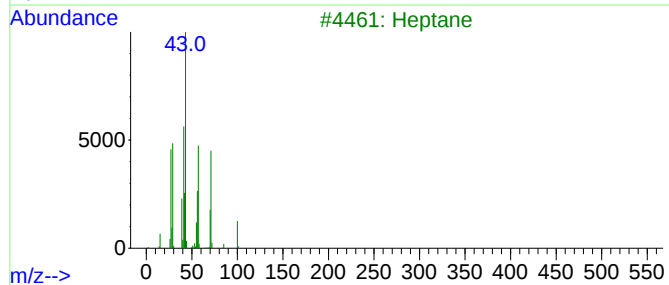
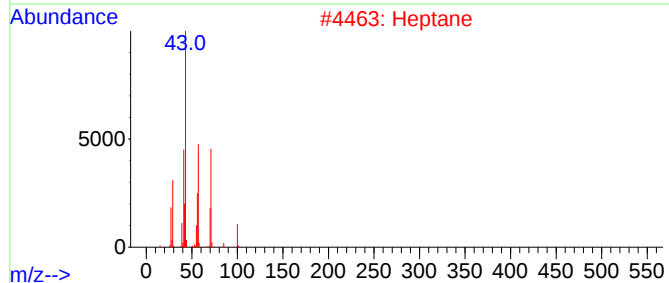
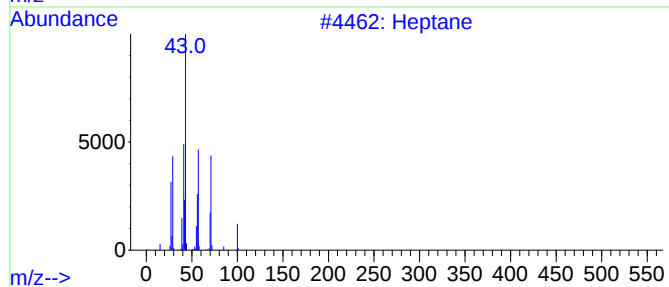
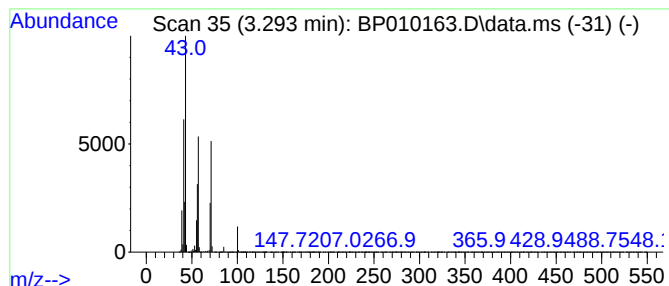
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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.
3.293	57.26 ng/ul	3473290	1,4-Dichlorobenzene-d4	7.922

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	83
5			Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	50



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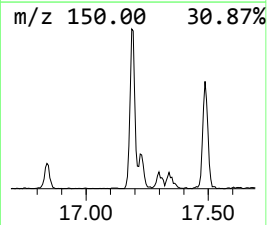
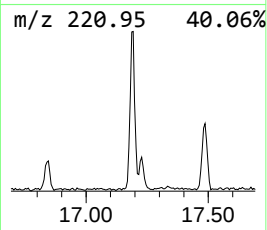
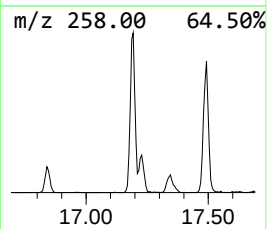
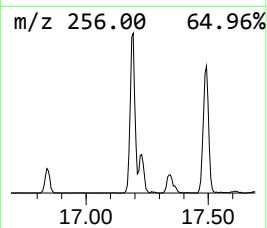
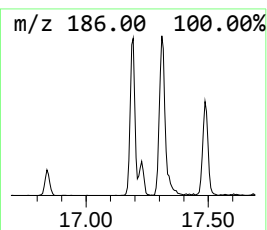
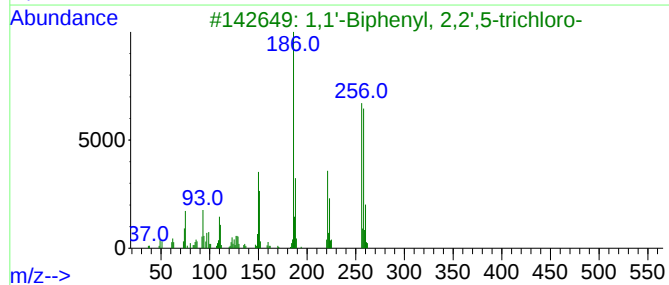
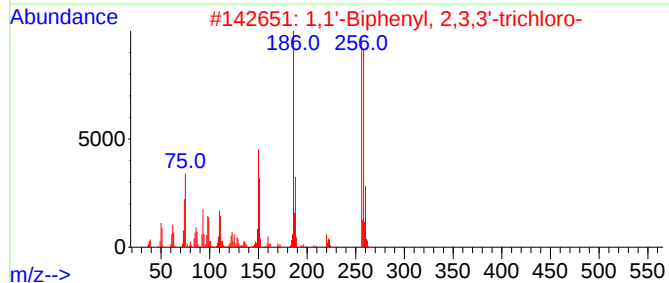
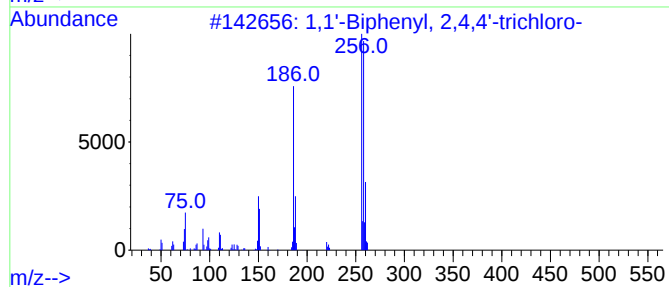
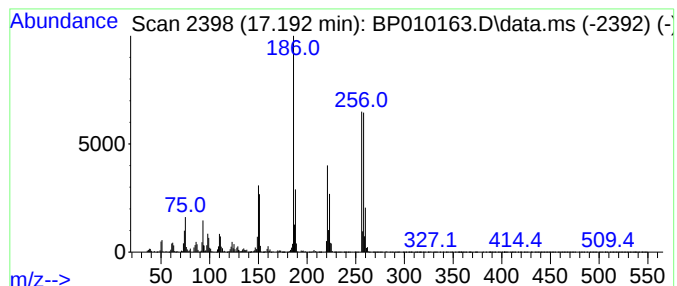
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.192	3.15 ng/ul	360998	Phenanthrene-d10	17.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
2	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	98		
3	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	97		
4	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	95		
5	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	95		



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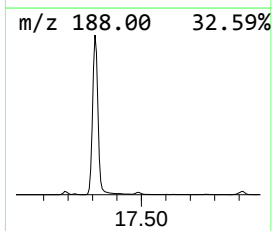
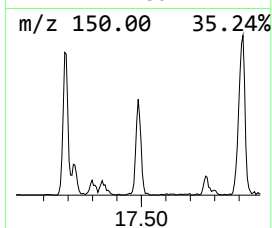
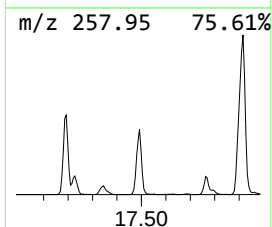
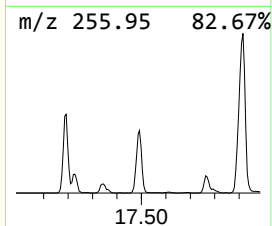
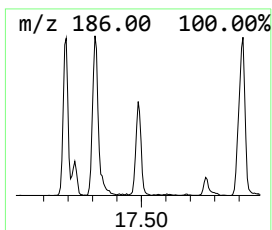
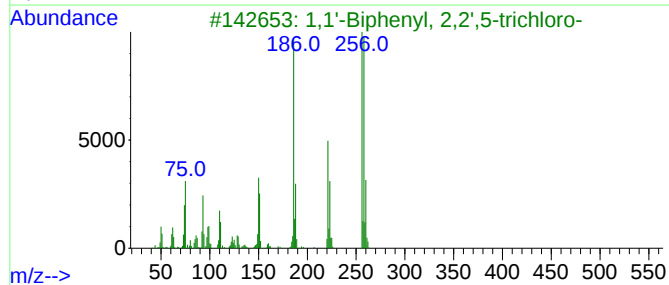
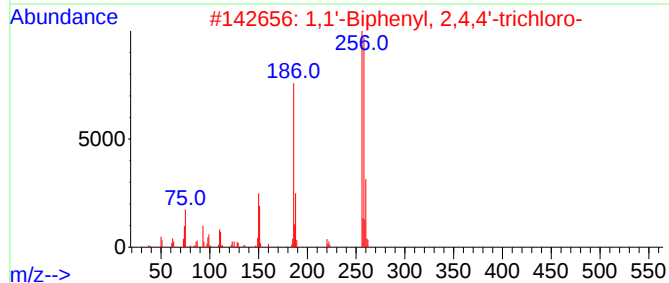
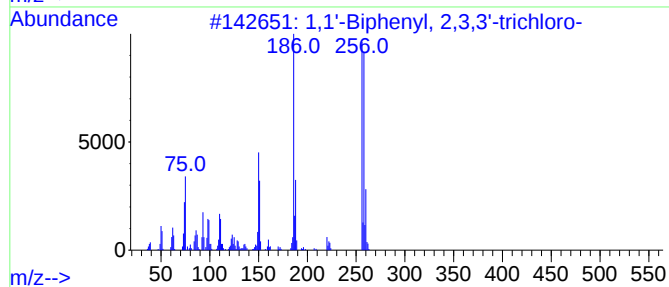
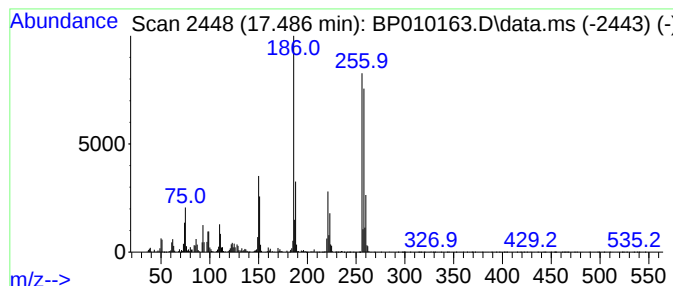
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TIC Library : C:\DATABASE\NIST20.L
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Peak Number 4 1,1'-Biphenyl, 2,3,3'-trich... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.486	2.26 ng/ul	259309	Phenanthrene-d10	17.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99		
2	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
3	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
4	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	97		
5	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	97		



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ClientSampleId :
ETNS5

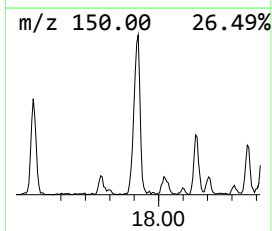
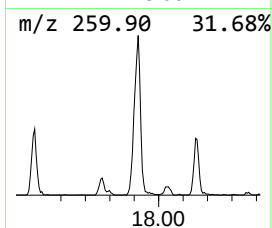
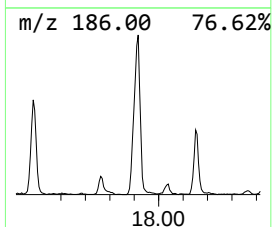
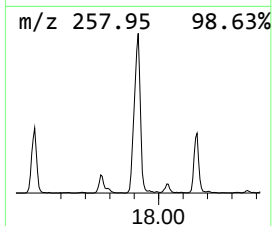
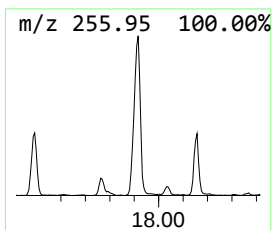
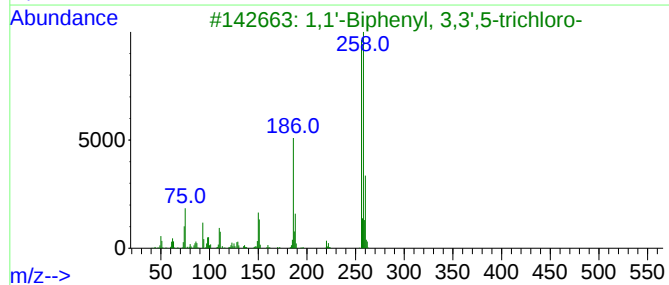
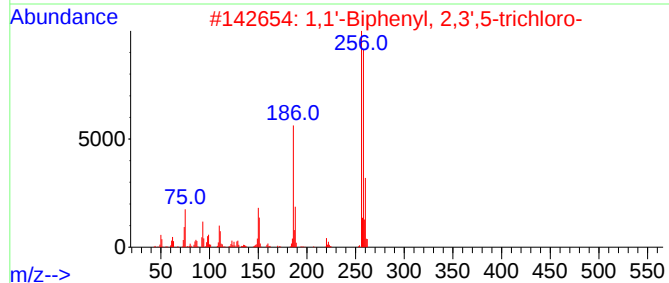
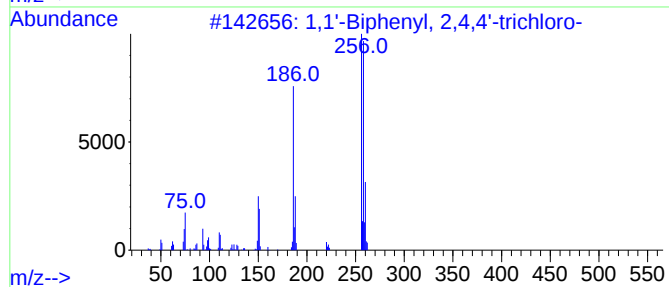
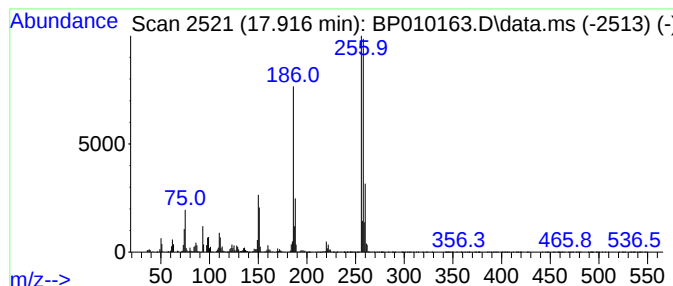
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.916	4.93 ng/ul	565393	Phenanthrene-d10	17.310

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99	
2	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99	
3	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99	
4	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99	
5	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
Data File : BP010163.D
Acq On : 29 Apr 2022 22:19
Operator : CG/JU
Sample : N2543-01
Misc : GCMS CONFIRMATION
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNS5

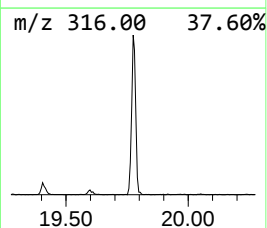
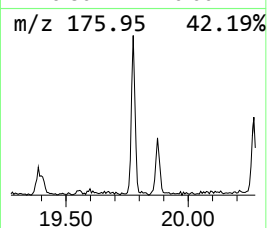
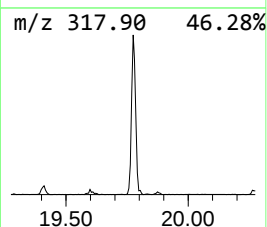
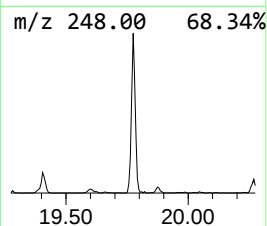
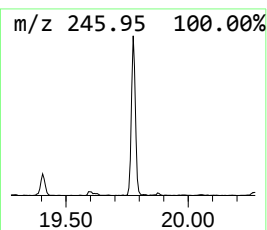
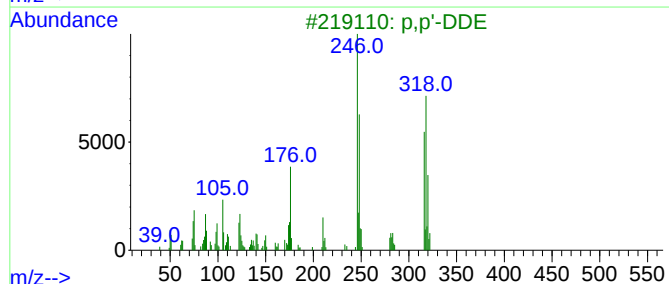
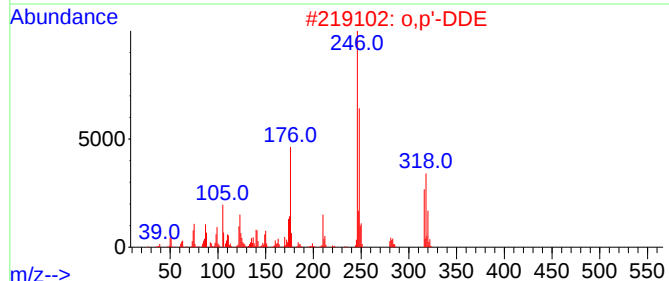
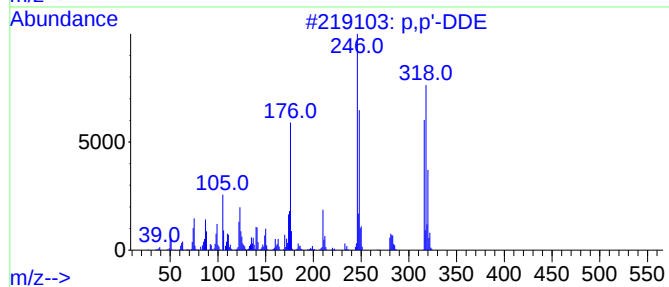
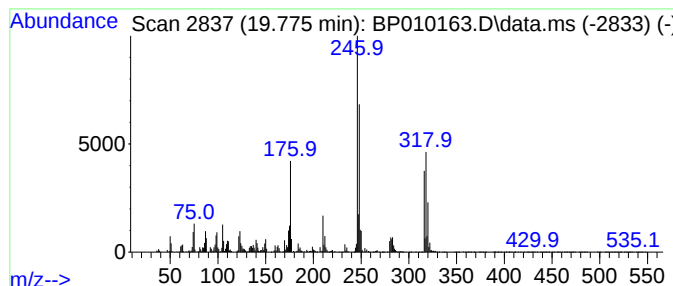
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 p,p'-DDE Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.775	2.78 ng/ul	268198	Chrysene-d12	21.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	p,p'-DDE			316	C14H8Cl4	000072-55-9	99
2	o,p'-DDE			316	C14H8Cl4	003424-82-6	99
3	p,p'-DDE			316	C14H8Cl4	000072-55-9	99
4	o,p'-DDE			316	C14H8Cl4	003424-82-6	95
5	p,p'-DDE			316	C14H8Cl4	000072-55-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
Data File : BP010163.D
Acq On : 29 Apr 2022 22:19
Operator : CG/JU
Sample : N2543-01
Misc : GCMS CONFIRMATION
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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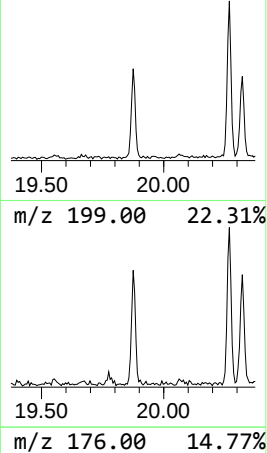
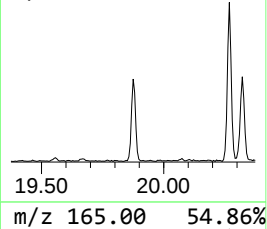
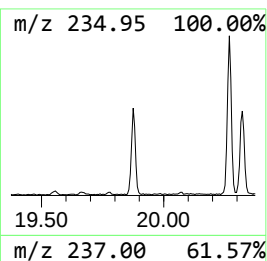
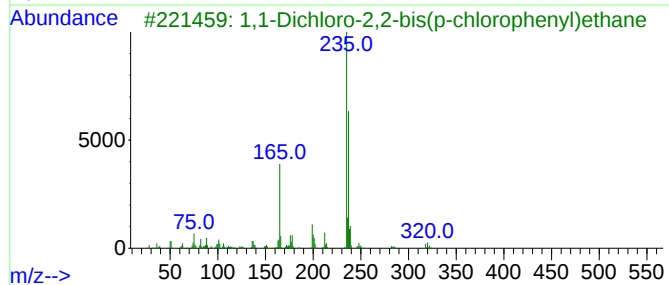
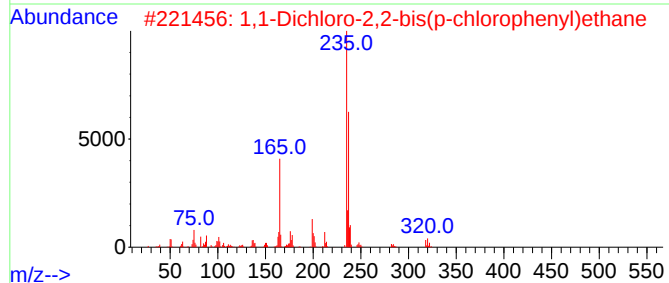
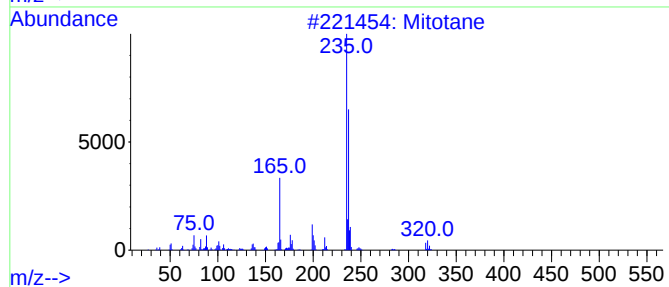
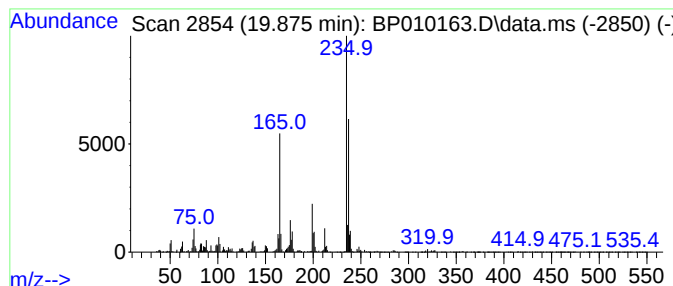
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Mitotane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.875	2.08 ng/ul	200640	Chrysene-d12	21.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mitotane	318	C14H10Cl4	000053-19-0	91
2			1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	90
3			1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	90
4			Mitotane	318	C14H10Cl4	000053-19-0	90
5			1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
Data File : BP010163.D
Acq On : 29 Apr 2022 22:19
Operator : CG/JU
Sample : N2543-01
Misc : GCMS CONFIRMATION
ALS Vial : 23 Sample Multiplier: 1

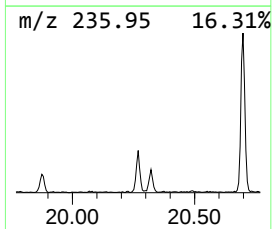
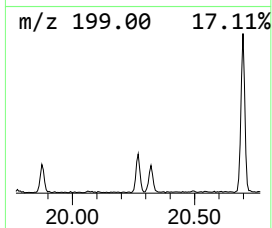
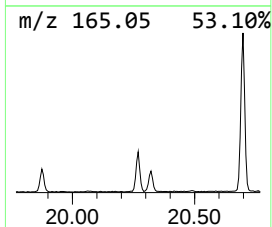
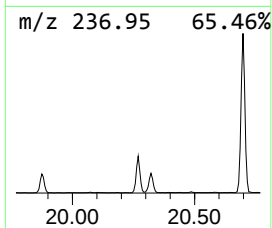
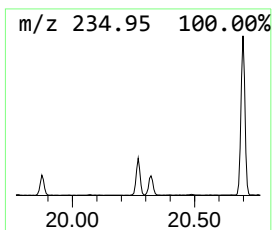
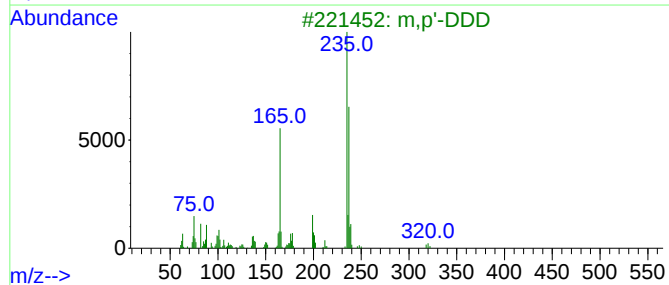
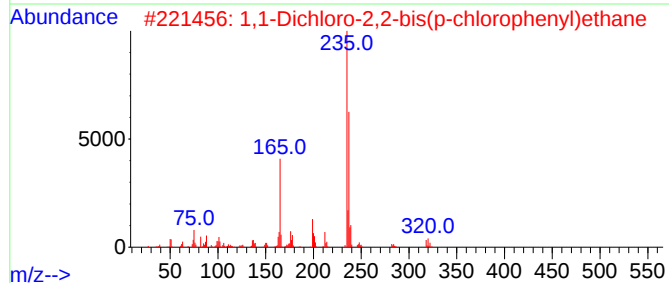
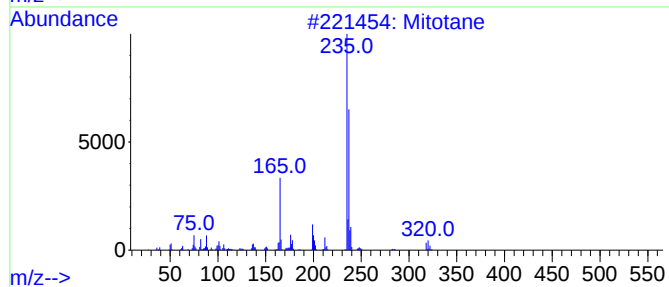
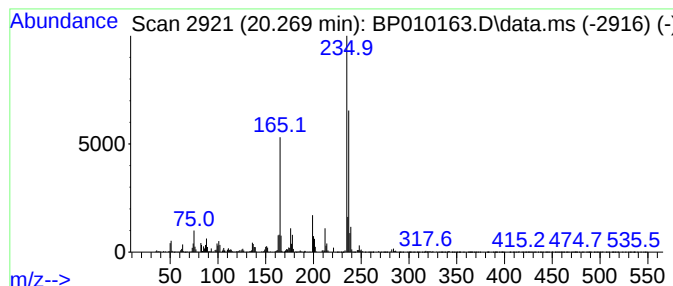
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 1,1-Dichloro-2,2-bis(p-chlo... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.269	3.32 ng/ul	320183	Chrysene-d12		21.392	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Mitotane		318	C14H10Cl4	000053-19-0	99
2	1,1-Dichloro-2,2-bis(p-chlorophe...		318	C14H10Cl4	000072-54-8	98
3	m,p'-DDD		318	C14H10Cl4	004329-12-8	91
4	p,p'-DDT		352	C14H9Cl5	000050-29-3	91
5	3-Butanone, 1,1-bis(4-chlorophen...		320	C18H18Cl2O	1000158-20-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
Data File : BP010163.D
Acq On : 29 Apr 2022 22:19
Operator : CG/JU
Sample : N2543-01
Misc : GCMS CONFIRMATION
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNS5

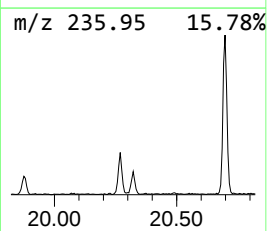
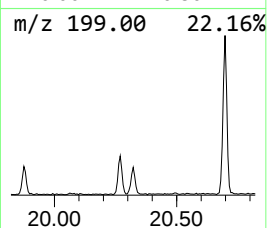
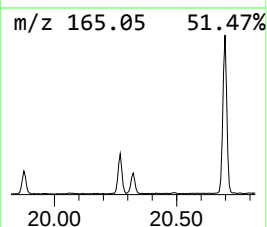
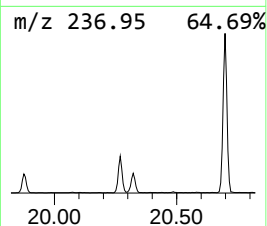
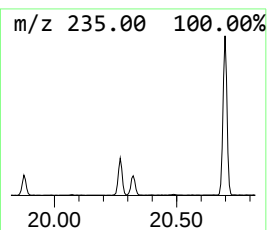
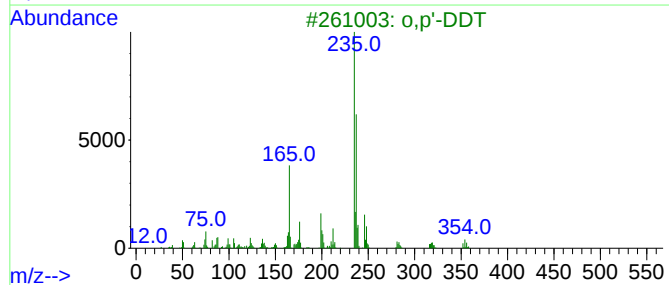
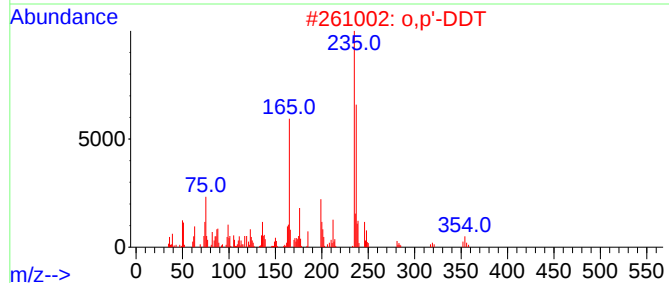
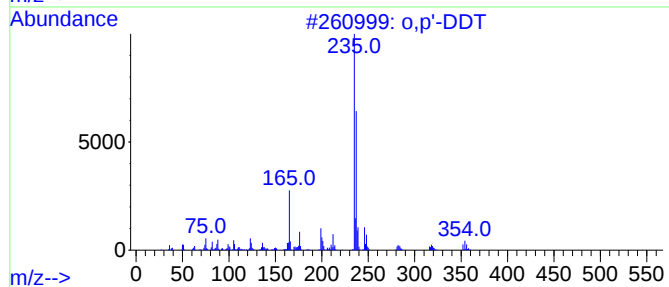
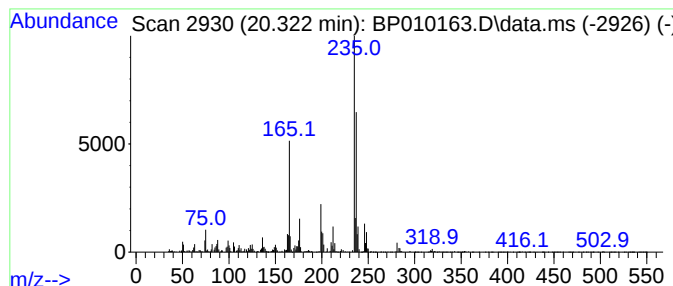
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 o,p'-DDT Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.322	2.10 ng/ul	202472	Chrysene-d12	21.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	o,p'-DDT			352	C14H9Cl5	000789-02-6	93
2	o,p'-DDT			352	C14H9Cl5	000789-02-6	91
3	o,p'-DDT			352	C14H9Cl5	000789-02-6	91
4	p,p'-DDT			352	C14H9Cl5	000050-29-3	91
5	Mitotane			318	C14H10Cl4	000053-19-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
Data File : BP010163.D
Acq On : 29 Apr 2022 22:19
Operator : CG/JU
Sample : N2543-01
Misc : GCMS CONFIRMATION
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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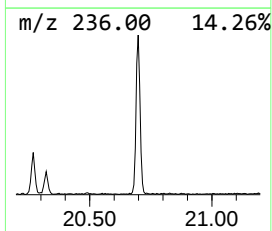
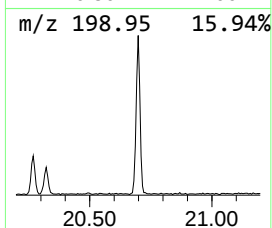
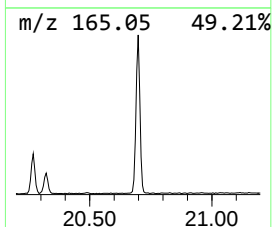
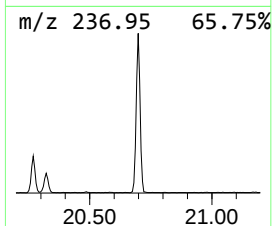
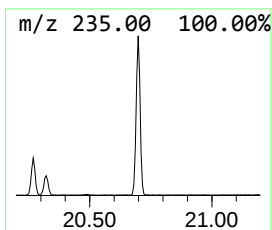
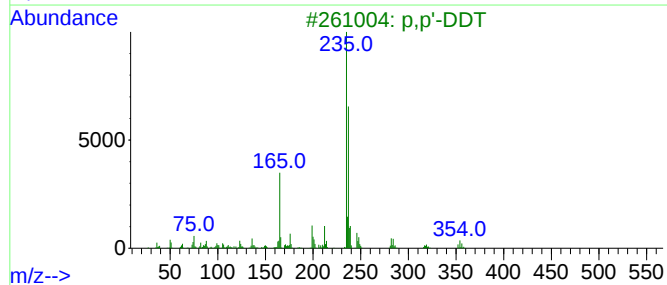
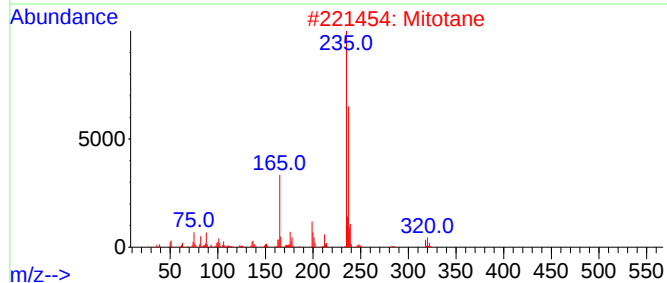
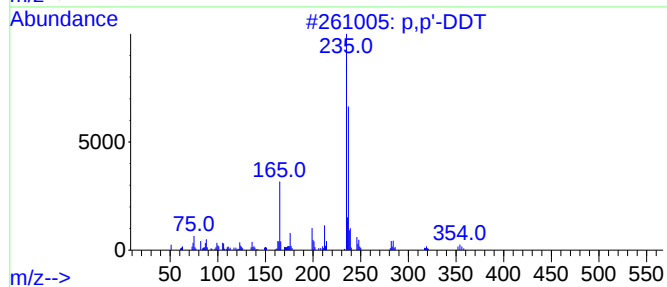
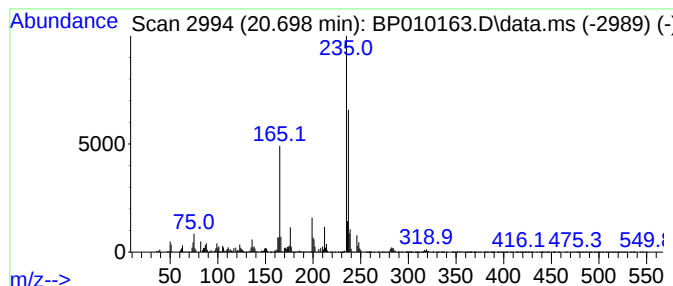
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 p,p'-DDT Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.698	14.52 ng/ul	1401910	Chrysene-d12	21.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	p,p'-DDT			352	C14H9Cl5	000050-29-3	99
2	Mitotane			318	C14H10Cl4	000053-19-0	98
3	p,p'-DDT			352	C14H9Cl5	000050-29-3	98
4	o,p'-DDT			352	C14H9Cl5	000789-02-6	97
5	1,1-Dichloro-2,2-bis(p-chlorophe...			318	C14H10Cl4	000072-54-8	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
 Data File : BP010163.D
 Acq On : 29 Apr 2022 22:19
 Operator : CG/JU
 Sample : N2543-01
 Misc : GCMS CONFIRMATION
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETNS5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.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
(DEL) Alkane: C...	3.117	3.4	ng/ul	205991	1	7.922	1213210	20.0
(DEL) Alkane: S...	3.293	57.3	ng/ul	3473290	1	7.922	1213210	20.0
1,1'-Biphenyl, ...	17.192	3.1	ng/ul	360998	4	17.310	2295690	20.0
1,1'-Biphenyl, ...	17.486	2.3	ng/ul	259309	4	17.310	2295690	20.0
1,1'-Biphenyl, ...	17.916	4.9	ng/ul	565393	4	17.310	2295690	20.0
p,p'-DDE	19.775	2.8	ng/ul	268198	5	21.392	1931320	20.0
Mitotane	19.875	2.1	ng/ul	200640	5	21.392	1931320	20.0
1,1-Dichloro-2,...	20.269	3.3	ng/ul	320183	5	21.392	1931320	20.0
o,p'-DDT	20.322	2.1	ng/ul	202472	5	21.392	1931320	20.0
p,p'-DDT	20.698	14.5	ng/ul	1401910	5	21.392	1931320	20.0