

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN052019\
 Data File : BN005807.D
 Acq On : 20 May 2019 19:44
 Operator : JU/SJ
 Sample : K2693-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BFH63

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.676	113	117	124	rVB	68399	92869	5.56%	0.418%
2	6.769	638	643	660	rVB	147653	358067	21.45%	1.610%
3	6.917	663	668	678	rBV	137078	228737	13.70%	1.029%
4	7.111	696	701	717	rVB	192195	349169	20.92%	1.570%
5	7.569	773	779	788	rBV	314440	540643	32.39%	2.431%
6	8.287	895	901	919	rVB	178529	376468	22.55%	1.693%
7	8.722	969	975	982	rBV	188069	340409	20.39%	1.531%
8	9.128	1042	1044	1046	rBV2	4614	3476	0.21%	0.016%
9	9.434	1091	1096	1109	rVB	154293	296687	17.77%	1.334%
10	9.987	1185	1190	1216	rBV2	147616	461455	27.64%	2.075%
11	10.328	1241	1248	1263	rVB	465494	817751	48.98%	3.678%
12	10.522	1276	1281	1291	rBV4	11963	32013	1.92%	0.144%
13	10.746	1317	1319	1322	rBV	2956	2441	0.15%	0.011%
14	12.028	1533	1537	1540	rBV4	5392	7334	0.44%	0.033%
15	12.234	1570	1572	1576	rBV3	3603	5153	0.31%	0.023%
16	12.457	1608	1610	1612	rBV2	2573	2995	0.18%	0.013%
17	12.622	1635	1638	1643	rBV	2001	3481	0.21%	0.016%
18	13.034	1705	1708	1713	rVB3	5597	8866	0.53%	0.040%
19	13.081	1713	1716	1718	rBV3	2403	2514	0.15%	0.011%
20	13.393	1760	1769	1774	rBV3	7334	16012	0.96%	0.072%
21	13.446	1774	1778	1781	rBV3	2940	4197	0.25%	0.019%
22	13.522	1788	1791	1796	rVB3	5371	8365	0.50%	0.038%
23	13.581	1796	1801	1803	rBV3	4954	8014	0.48%	0.036%
24	13.622	1803	1808	1813	rVV	506730	765046	45.83%	3.441%
25	13.675	1813	1817	1825	rVB	127905	206105	12.35%	0.927%
26	13.769	1832	1833	1838	rVB4	3591	3957	0.24%	0.018%
27	13.898	1847	1855	1864	rBV	616888	959669	57.49%	4.316%
28	14.116	1889	1892	1898	rBV3	7326	11369	0.68%	0.051%
29	14.204	1901	1907	1916	rVV2	795282	1220538	73.11%	5.489%
30	14.316	1924	1926	1930	rBV4	3303	3163	0.19%	0.014%
31	14.481	1948	1954	1965	rBV2	72992	217465	13.03%	0.978%
32	14.622	1973	1978	1987	rVB8	20646	44549	2.67%	0.200%
33	14.751	1995	2000	2002	rBV4	3407	5287	0.32%	0.024%
34	14.840	2010	2015	2019	rBV4	5385	5770	0.35%	0.026%

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35	14.893	2021	2024	2027	rBV4	4820	6477	0.39%	0.029%
36	14.934	2029	2031	2034	rVB4	2443	2435	0.15%	0.011%
37	14.975	2034	2038	2041	rBV4	9329	15292	0.92%	0.069%
38	15.081	2053	2056	2061	rVB3	13166	17889	1.07%	0.080%
39	15.210	2071	2078	2085	rBV	838660	1254530	75.15%	5.642%
40	15.263	2085	2087	2089	rVV3	6635	6539	0.39%	0.029%
41	15.322	2093	2097	2101	rVB3	4308	6466	0.39%	0.029%
42	15.381	2101	2107	2114	rBV	77455	156889	9.40%	0.706%
43	15.451	2116	2119	2125	rBV4	8775	13982	0.84%	0.063%
44	15.575	2136	2140	2141	rBV3	5335	5518	0.33%	0.025%
45	15.640	2148	2151	2155	rBV6	4337	7120	0.43%	0.032%
46	15.722	2157	2165	2166	rBV8	8568	16156	0.97%	0.073%
47	15.740	2166	2168	2171	rVB4	5784	5663	0.34%	0.025%
48	15.804	2177	2179	2180	rBV2	3039	2649	0.16%	0.012%
49	15.945	2199	2203	2206	rBV4	20549	31784	1.90%	0.143%
50	16.040	2217	2219	2221	rBV3	4024	4800	0.29%	0.022%
51	16.116	2231	2232	2235	rBV3	8306	8104	0.49%	0.036%
52	16.251	2253	2255	2256	rBV2	6419	3761	0.23%	0.017%
53	16.516	2298	2300	2303	rBV4	8679	11358	0.68%	0.051%
54	16.545	2303	2305	2308	rVV4	9469	10003	0.60%	0.045%
55	16.740	2335	2338	2342	rBV5	18610	31839	1.91%	0.143%
56	16.963	2371	2376	2382	rBV2	1178103	1669421	100.00%	7.508%
57	17.069	2388	2394	2404	rBV	913699	1405819	84.21%	6.322%
58	17.887	2529	2533	2542	rBV3	200269	369883	22.16%	1.663%
59	19.186	2751	2754	2763	rBV5	154525	315473	18.90%	1.419%
60	19.386	2783	2788	2794	rBV	1108760	1637432	98.08%	7.364%
61	19.769	2850	2853	2856	rVB	265930	280483	16.80%	1.261%
62	19.910	2873	2877	2882	rBV	731827	837169	50.15%	3.765%
63	20.192	2922	2925	2930	rVB	541925	621360	37.22%	2.794%
64	20.351	2949	2952	2957	rVB2	309653	415284	24.88%	1.868%
65	20.516	2974	2980	2986	rVB	500475	868684	52.04%	3.907%
66	20.663	3002	3005	3009	rVB	467029	496686	29.75%	2.234%
67	20.933	3048	3051	3057	rBV2	545304	718528	43.04%	3.231%
68	21.210	3094	3098	3100	rBV	1206844	1640917	98.29%	7.379%
69	21.545	3152	3155	3160	rVB	336113	429897	25.75%	1.933%
70	23.292	3447	3452	3459	rVB3	699223	1500049	89.85%	6.746%

LSC Area Percent Report

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Sampling : 1 Min Area: 0.1 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M
Title : SVOA CALIBRATION

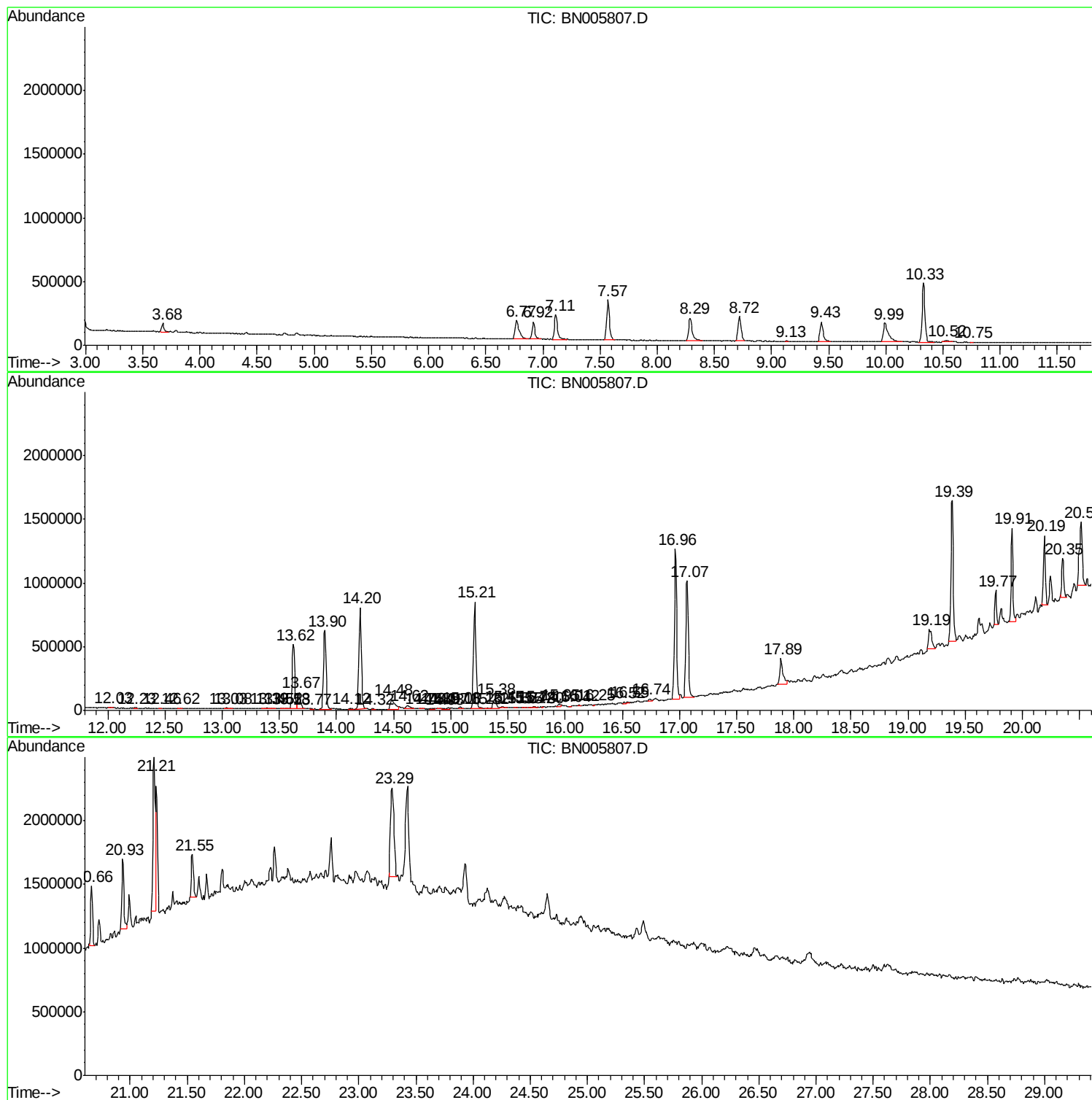
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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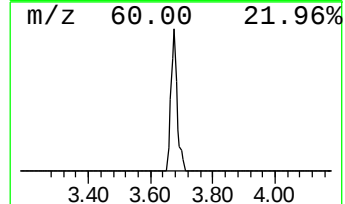
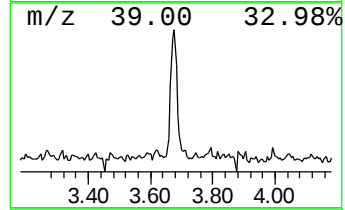
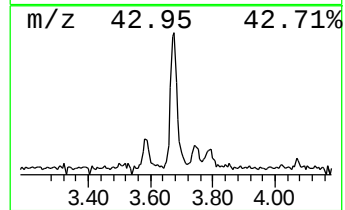
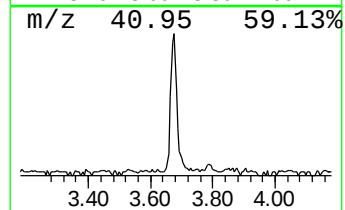
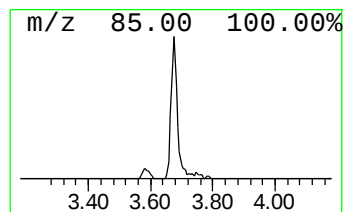
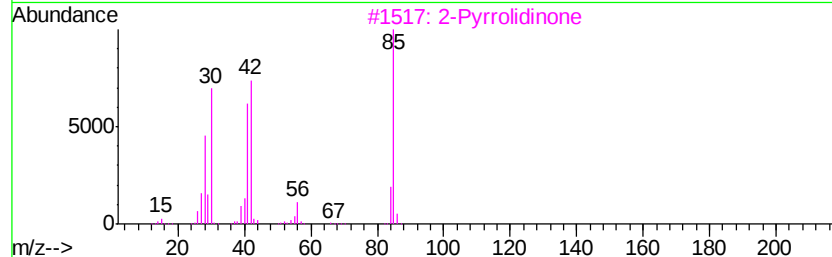
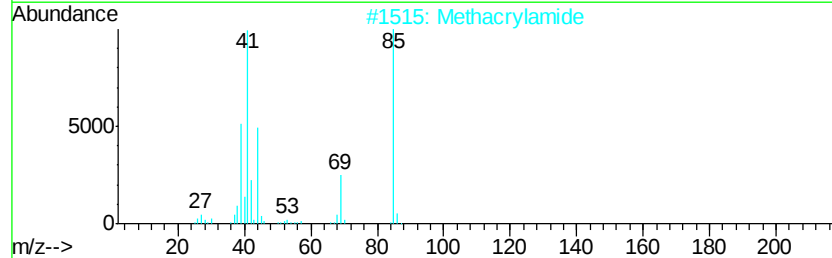
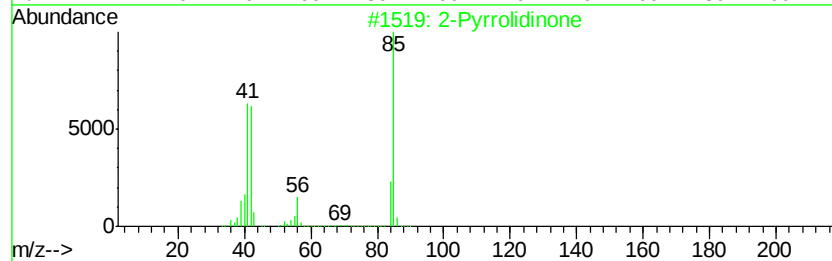
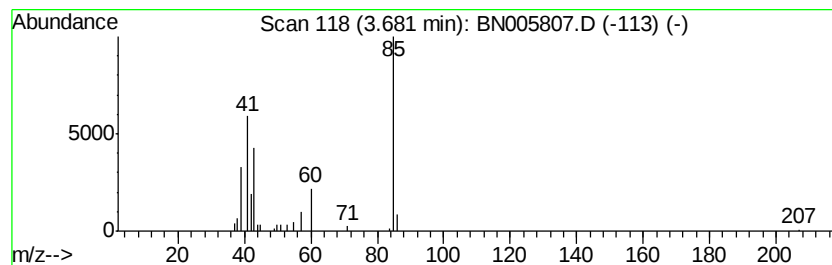
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.68	3.44 ng/ul	92869	1,4-Dichlorobenzene-d4	7.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pyrrolidinone			85	C4H7NO	000616-45-5	40
2	Methacrylamide			85	C4H7NO	000079-39-0	38
3	2-Pyrrolidinone			85	C4H7NO	000616-45-5	28
4	Methacrylamide			85	C4H7NO	000079-39-0	10
5	2-Pyrrolidinone			85	C4H7NO	000616-45-5	9



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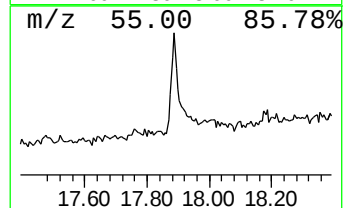
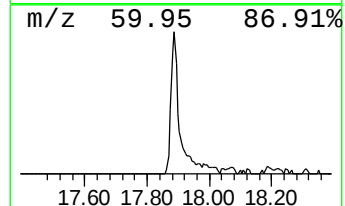
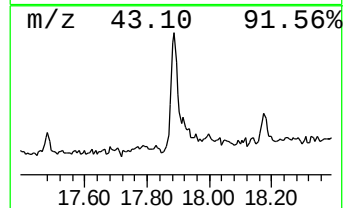
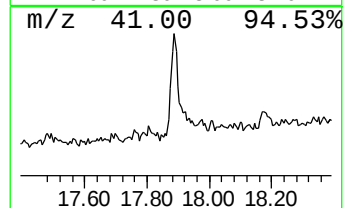
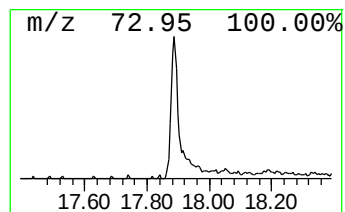
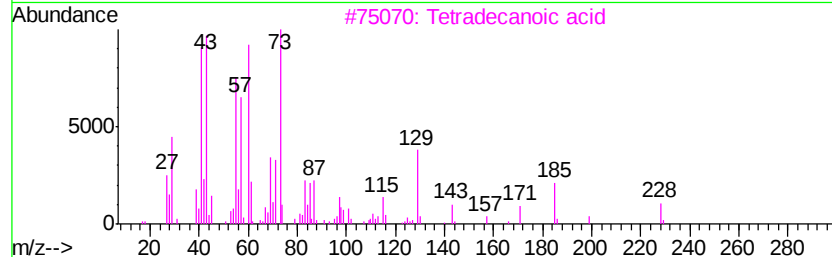
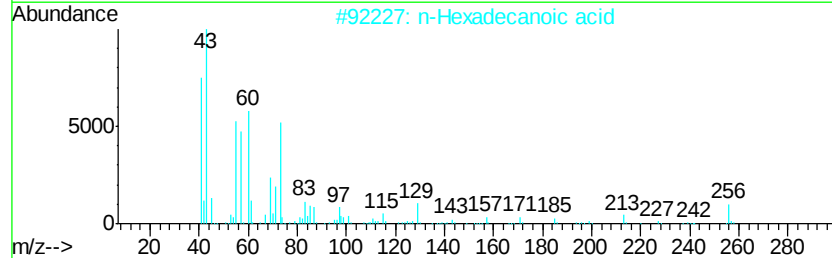
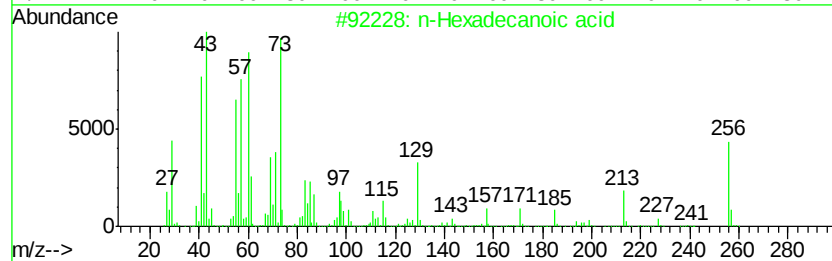
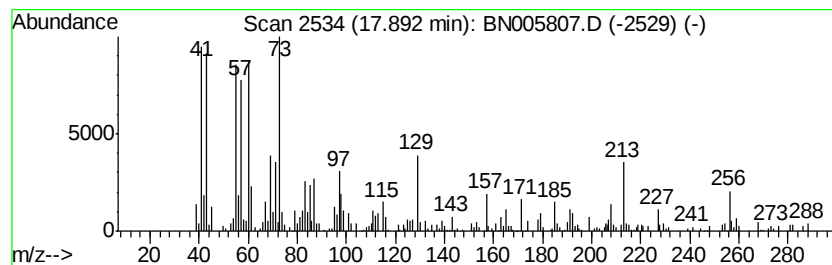
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.89	4.43 ng/ul	369883	Phenanthrene-d10	16.96		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97	
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94	
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	89	
4	Octadecanoic acid	284	C18H36O2	000057-11-4	81	
5	Undecanoic acid	186	C11H22O2	000112-37-8	78	



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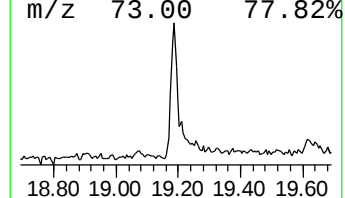
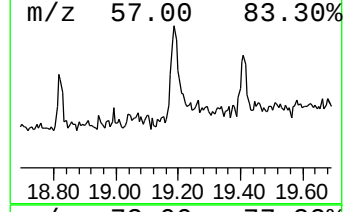
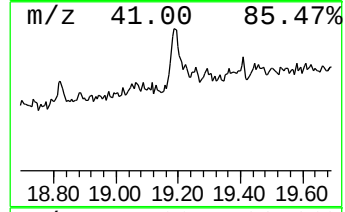
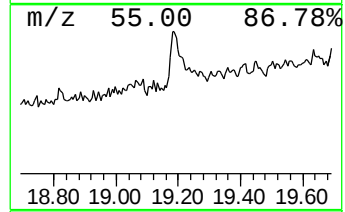
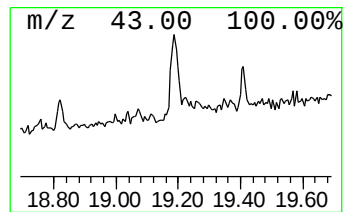
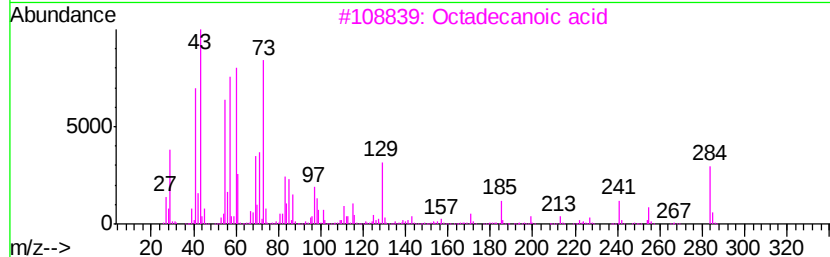
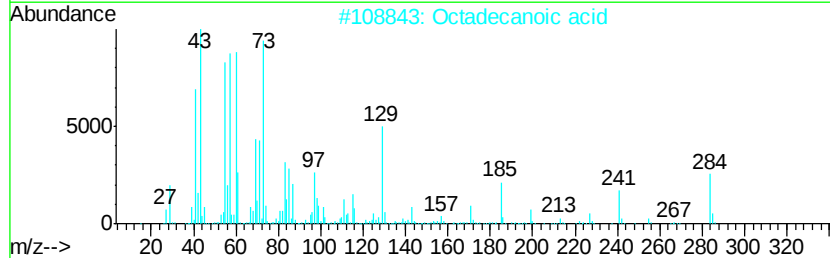
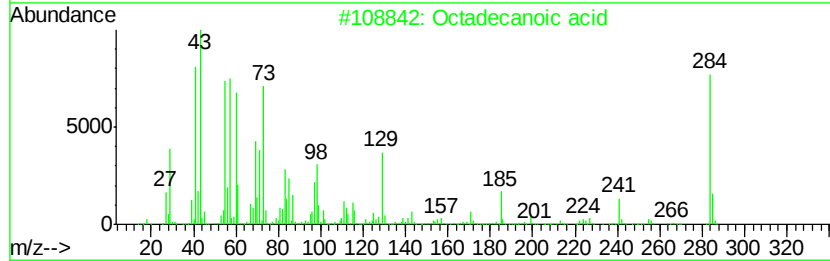
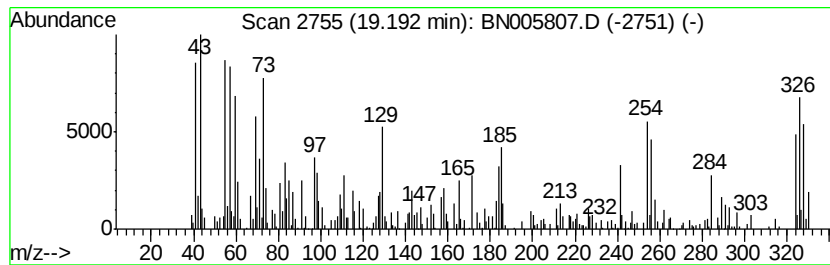
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.19	3.85 ng/ul	315473	Chrysene-d12	21.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	84
2		Octadecanoic acid	284	C18H36O2	000057-11-4	84
3		Octadecanoic acid	284	C18H36O2	000057-11-4	81
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	80
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	42



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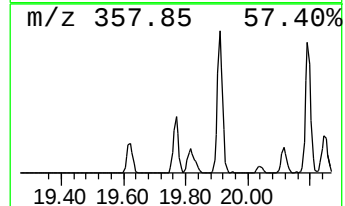
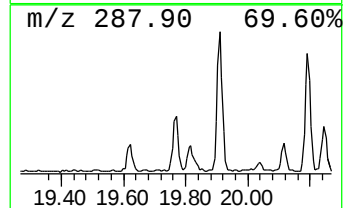
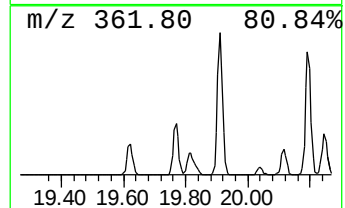
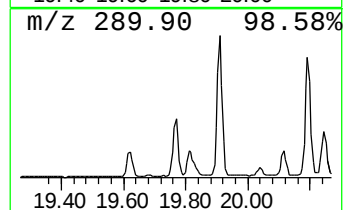
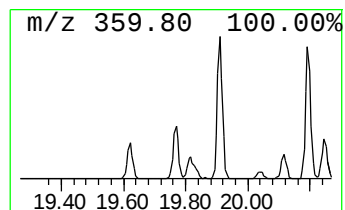
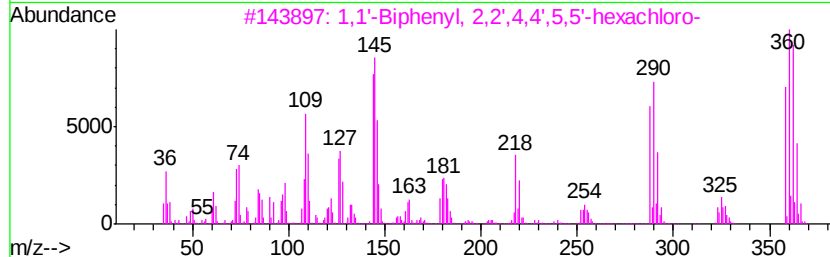
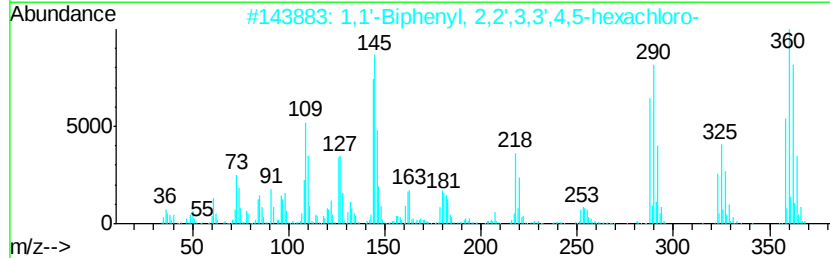
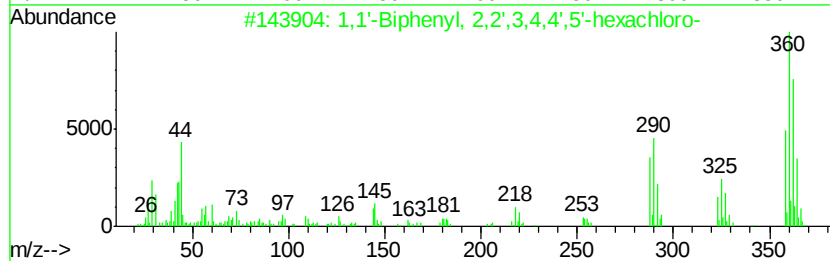
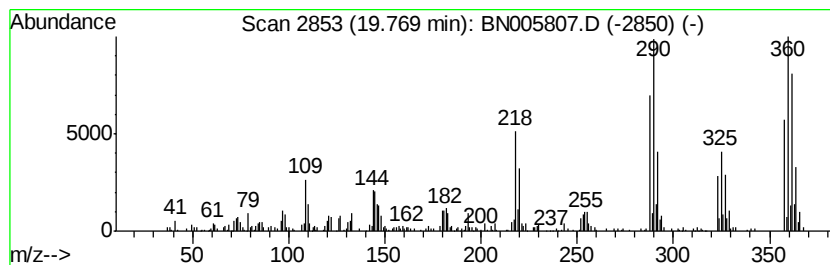
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.77	3.42 ng/ul	280483	Chrysene-d12	21.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,4,4',5'-he...	358	C12H4Cl6	035065-28-2	98
2		1,1'-Biphenyl, 2,2',3,3',4,5-hex...	358	C12H4Cl6	055215-18-4	97
3		1,1'-Biphenyl, 2,2',4,4',5,5'-he...	358	C12H4Cl6	035065-27-1	94
4		1,1'-Biphenyl, 2,2',3,4,4',5',6-he...	358	C12H4Cl6	038380-04-0	94
5		1,1'-Biphenyl, 2,2',4,4',5,5'-he...	358	C12H4Cl6	035065-27-1	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN052019\
Data File : BN005807.D
Acq On : 20 May 2019 19:44
Operator : JU/SJ
Sample : K2693-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BFH63

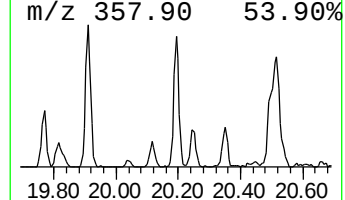
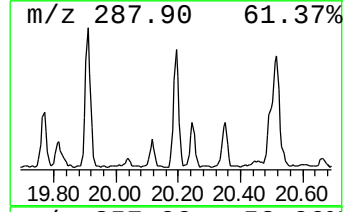
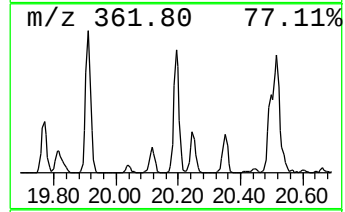
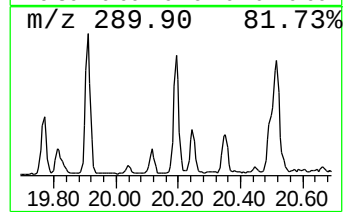
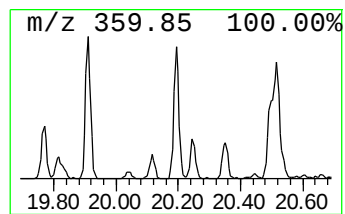
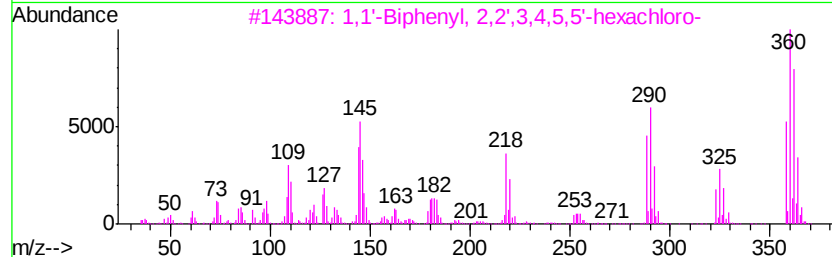
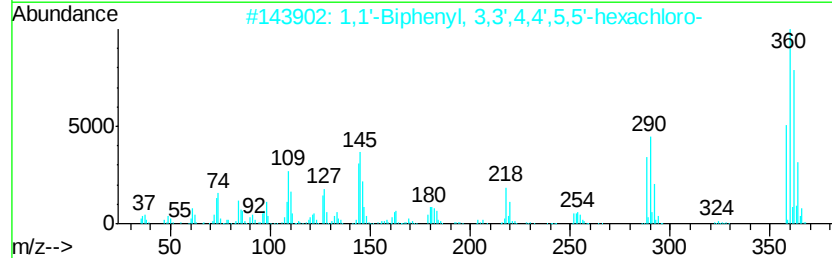
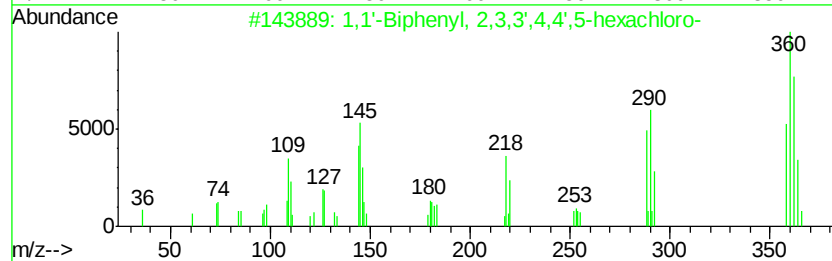
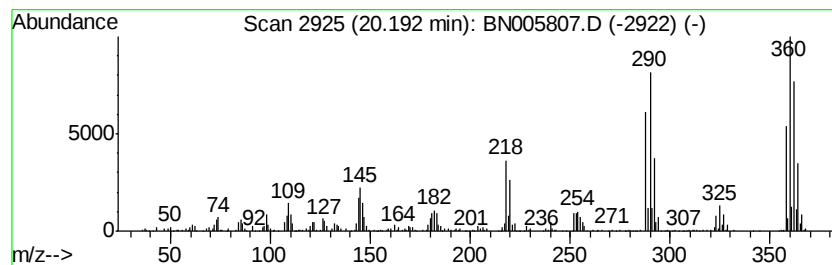
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.19	7.57 ng/ul	621360	Chrysene-d12	21.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99
2		1,1'-Biphenyl, 3,3',4,4',5,5'-he...	358	C12H4Cl6	032774-16-6	99
3		1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	96
4		1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	96
5		1,1'-Biphenyl, 2,2',3,3',5,6'-he...	358	C12H4Cl6	052744-13-5	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN052019\
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Operator : JU/SJ
Sample : K2693-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

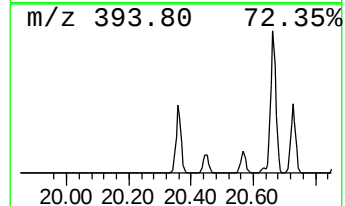
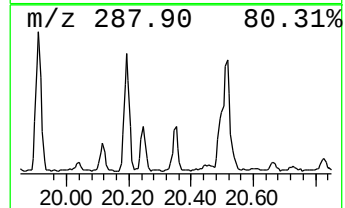
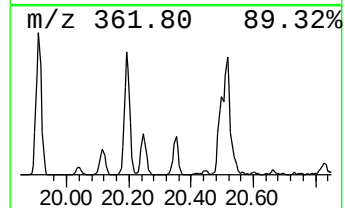
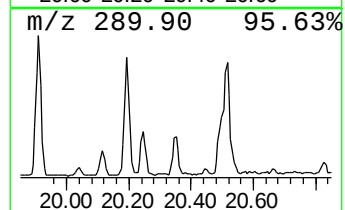
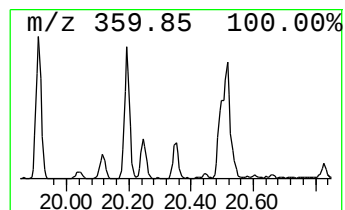
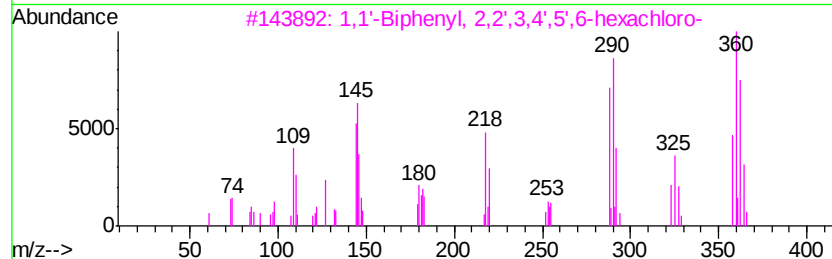
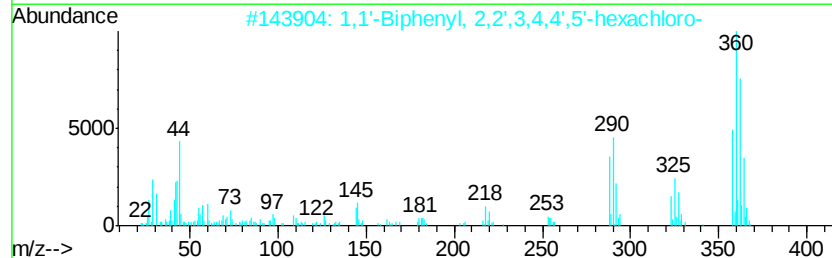
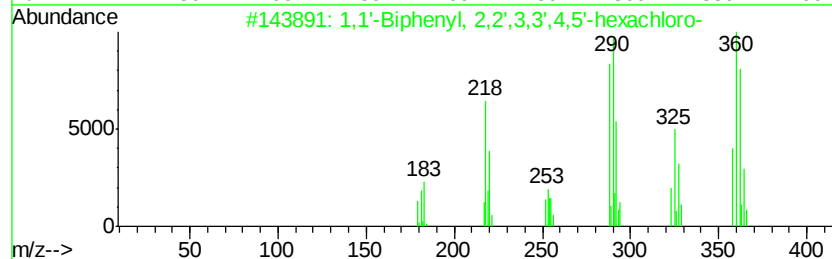
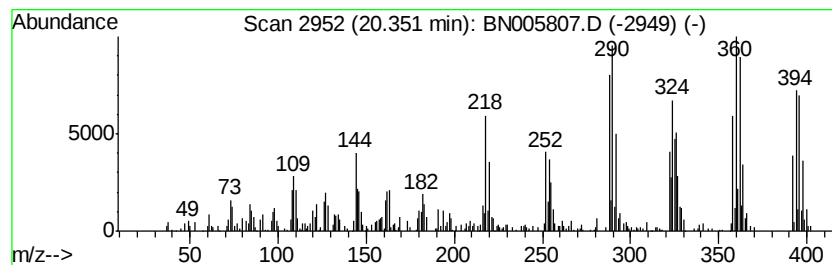
Instrument :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.35	5.06 ng/ul	415284	Chrysene-d12	21.21		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl,	2,2',3,3',4,5'-he...	358	C12H4Cl6	052663-66-8	99
2	1,1'-Biphenyl,	2,2',3,4,4',5'-he...	358	C12H4Cl6	035065-28-2	96
3	1,1'-Biphenyl,	2,2',3,4',5',6-he...	358	C12H4Cl6	038380-04-0	94
4	1,1'-Biphenyl,	2,2',3,4,5,6'-Hex...	358	C12H4Cl6	068194-15-0	94
5	1,1'-Biphenyl,	2,2',3,3',4,5-hex...	358	C12H4Cl6	055215-18-4	94



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Sample : K2693-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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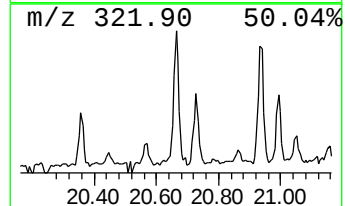
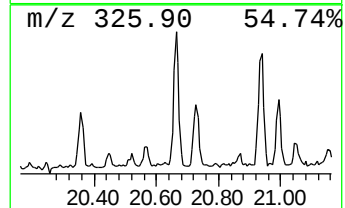
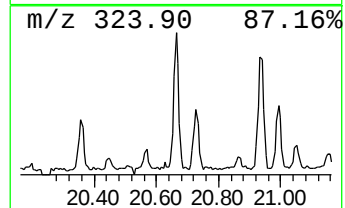
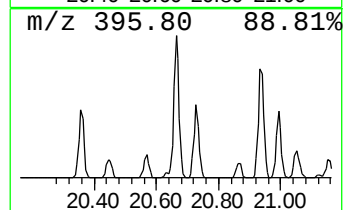
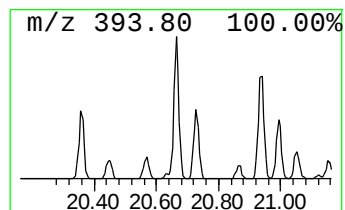
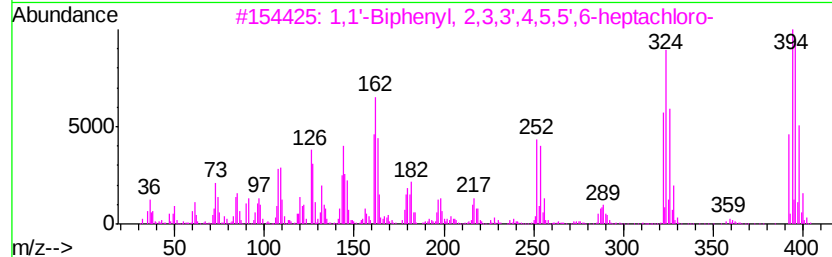
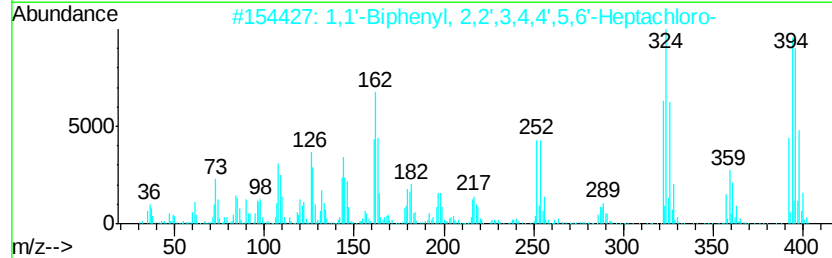
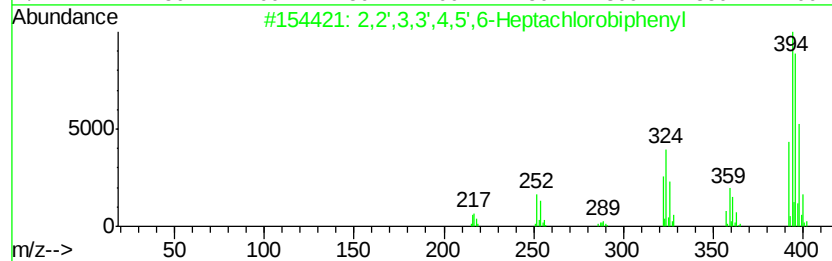
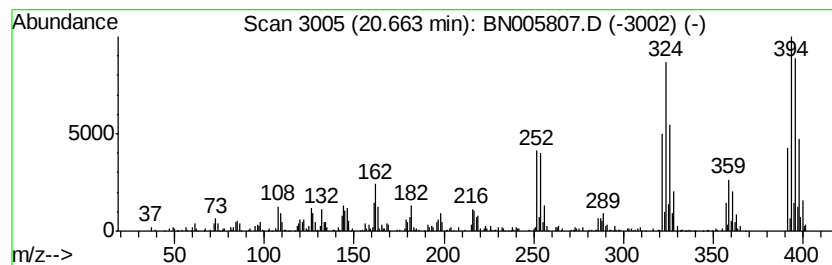
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.66	6.05 ng/ul	496686	Chrysene-d12	21.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2',3,3',4,5',6-Heptachlorobiph...	392	C12H3Cl7	040186-70-7	99
2		1,1'-Biphenyl, 2,2',3,4,4',5,6'-...	392	C12H3Cl7	060145-23-5	99
3		1,1'-Biphenyl, 2,3,3',4,5,5',6-h...	392	C12H3Cl7	074472-51-8	99
4		1,1'-Biphenyl, 2,2',3,4,4',5,5'-...	392	C12H3Cl7	035065-29-3	99
5		1,1'-Biphenyl, 2,2',3,4,5,6,6'-H...	392	C12H3Cl7	074472-49-4	96



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Sample : K2693-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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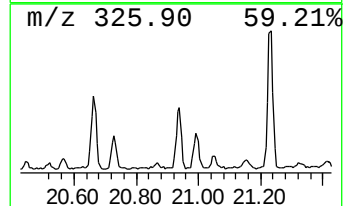
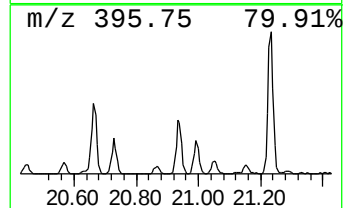
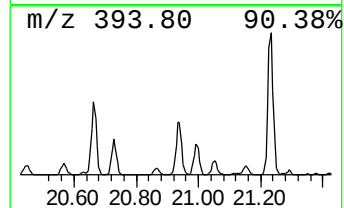
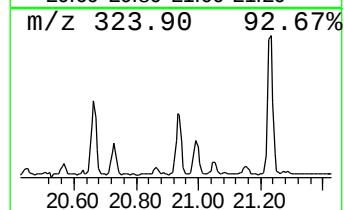
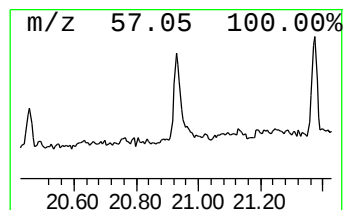
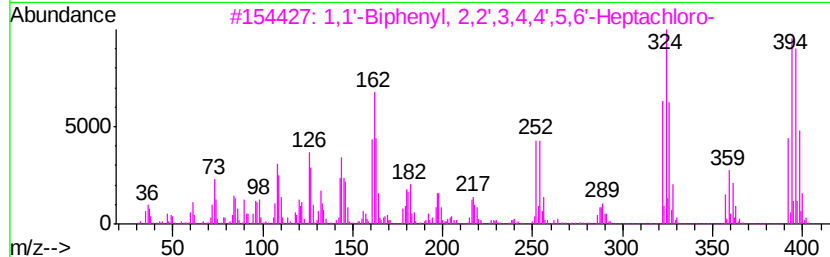
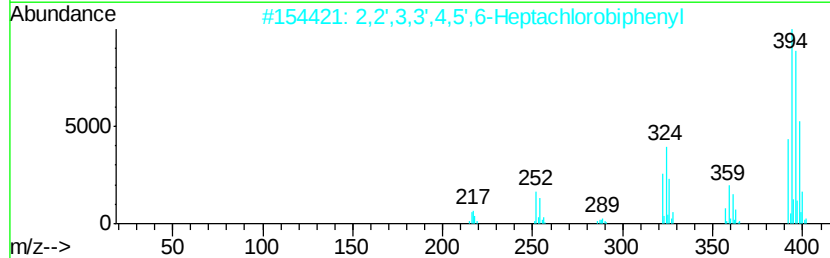
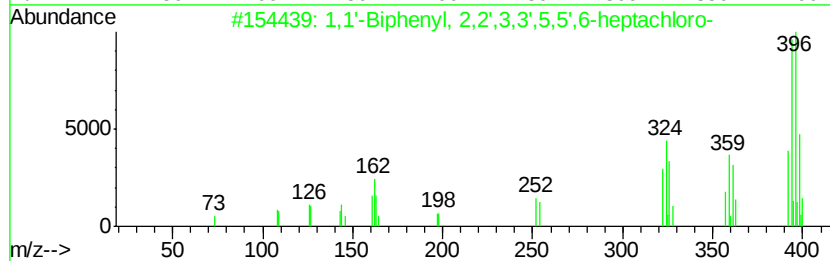
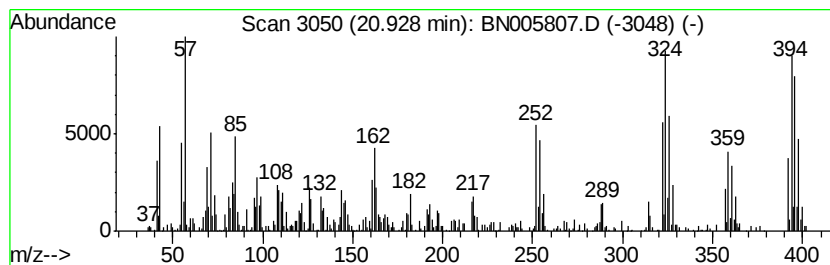
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.93	8.76 ng/ul	718528	Chrysene-d12	21.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,3',5,5',6-...	392	C12H3Cl7	052663-67-9	99
2		2,2',3,3',4,5',6-Heptachlorobiph...	392	C12H3Cl7	040186-70-7	99
3		1,1'-Biphenyl, 2,2',3,4,4',5,6'-...	392	C12H3Cl7	060145-23-5	99
4		1,1'-Biphenyl, 2,2',3,4,5,5',6-h...	392	C12H3Cl7	052712-05-7	99
5		1,1'-Biphenyl, 2,3,3',4,5,5',6-h...	392	C12H3Cl7	074472-51-8	99



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Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	3.68	3.4	ng/ul	92869	1	7.57	540643	20.0
n-Hexadecanoic acid	17.89	4.4	ng/ul	369883	4	16.96	1669420	20.0
Octadecanoic acid	19.19	3.9	ng/ul	315473	5	21.21	1640920	20.0
1,1'-Biphenyl, 2,...	19.77	3.4	ng/ul	280483	5	21.21	1640920	20.0
1,1'-Biphenyl, 2,...	20.19	7.6	ng/ul	621360	5	21.21	1640920	20.0
1,1'-Biphenyl, 2,...	20.35	5.1	ng/ul	415284	5	21.21	1640920	20.0
2,2',3,3',4,5',6-...	20.66	6.0	ng/ul	496686	5	21.21	1640920	20.0
1,1'-Biphenyl, 2,...	20.93	8.8	ng/ul	718528	5	21.21	1640920	20.0