

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060322\
 Data File : BG053811.D
 Acq On : 3 Jun 2022 13:40
 Operator : CG/JU
 Sample : N2984-17
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P021-SS001-0612-01

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % 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_G\Methods\8270-BG060222.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG053811.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.656	377	386	398	rBV	417586	711721	37.21%	6.877%
2	7.243	648	656	665	rBV	429196	771447	40.33%	7.454%
3	8.095	794	801	807	rBV	90201	159109	8.32%	1.537%
4	9.252	988	998	1006	rBV	256185	486406	25.43%	4.700%
5	10.903	1271	1279	1289	rVB	128681	239915	12.54%	2.318%
6	13.342	1686	1694	1702	rVB	803669	1268529	66.32%	12.257%
7	14.711	1921	1927	1934	rBV	214364	351347	18.37%	3.395%
8	15.269	2010	2022	2034	rBV2	7518	21362	1.12%	0.206%
9	16.197	2173	2180	2190	rBV	567753	941408	49.22%	9.096%
10	17.454	2387	2394	2400	rBV	269969	430517	22.51%	4.160%
11	18.342	2541	2545	2553	rBV2	32035	50448	2.64%	0.487%
12	19.505	2739	2743	2749	rVB	19957	30915	1.62%	0.299%
13	19.646	2764	2767	2772	rVB2	29488	37211	1.95%	0.360%
14	19.863	2798	2804	2809	rVB2	18101	40161	2.10%	0.388%
15	20.063	2832	2838	2846	rBV	1373053	1912747	100.00%	18.482%
16	20.210	2860	2863	2868	rBV	40799	55100	2.88%	0.532%
17	20.257	2868	2871	2877	rVB7	16290	26786	1.40%	0.259%
18	20.351	2882	2887	2900	rBV2	117605	187270	9.79%	1.809%
19	20.551	2917	2921	2924	rBV4	22361	29075	1.52%	0.281%
20	20.627	2929	2934	2939	rVV2	146728	186769	9.76%	1.805%
21	20.680	2939	2943	2948	rVV4	33186	43543	2.28%	0.421%
22	20.786	2956	2961	2968	rBV5	48533	96063	5.02%	0.928%
23	20.880	2973	2977	2981	rVV3	25036	40998	2.14%	0.396%
24	20.950	2981	2989	2996	rVV2	80108	194531	10.17%	1.880%
25	21.009	2996	2999	3004	rVB6	16165	20489	1.07%	0.198%
26	21.115	3012	3017	3024	rVB2	88478	130579	6.83%	1.262%
27	21.179	3024	3028	3032	rBV3	36842	54860	2.87%	0.530%
28	21.379	3057	3062	3066	rBV2	89368	147705	7.72%	1.427%
29	21.414	3066	3068	3075	rVB3	66040	92311	4.83%	0.892%
30	21.485	3076	3080	3088	rVB4	39601	63083	3.30%	0.610%
31	21.556	3088	3092	3095	rBV5	16186	25857	1.35%	0.250%
32	21.732	3116	3122	3126	rBV2	278810	501120	26.20%	4.842%
33	21.773	3126	3129	3135	rVB2	130108	198315	10.37%	1.916%
34	21.949	3154	3159	3163	rBV2	19571	34113	1.78%	0.330%
35	22.184	3195	3199	3208	rVB4	51020	102702	5.37%	0.992%
36	22.266	3210	3213	3220	rVB8	22701	41157	2.15%	0.398%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_G\Methods\8270-BG060222.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	22.366	3226	3230	3236	rVB7	25877	38362	2.01%	0.371%
38	22.572	3261	3265	3273	rVB6	20972	41387	2.16%	0.400%
39	24.111	3522	3527	3532	rBV2	16746	34877	1.82%	0.337%
40	25.004	3668	3679	3689	rBV2	151768	460676	24.08%	4.451%
41	26.262	3885	3893	3900	rBV5	14536	48399	2.53%	0.468%

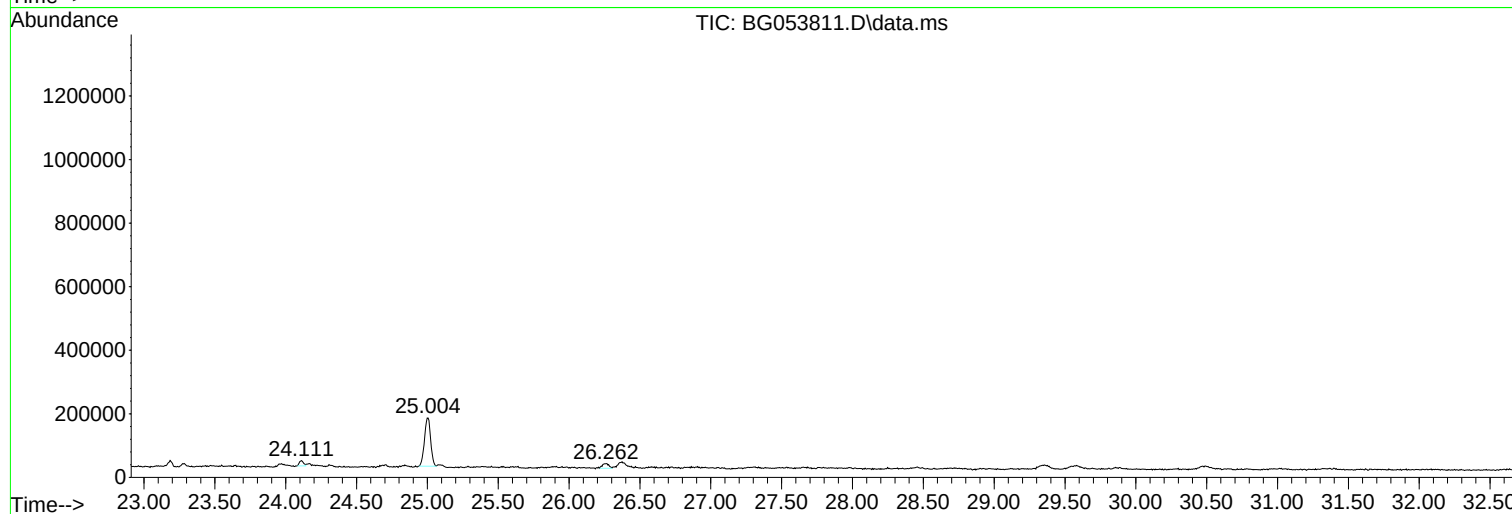
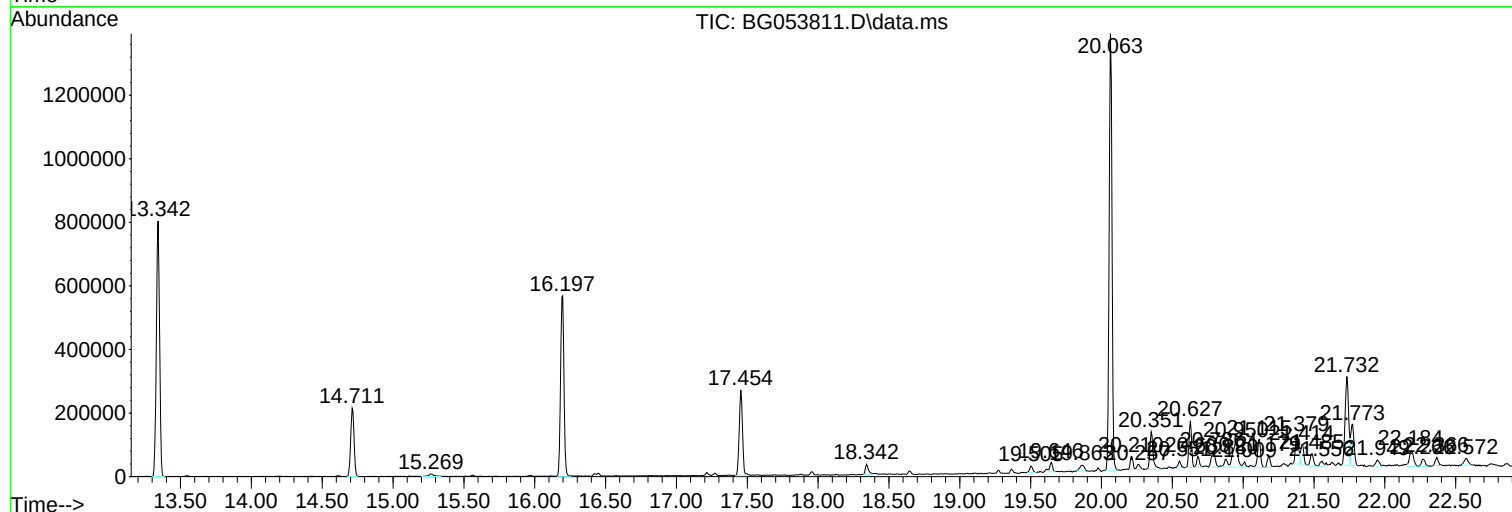
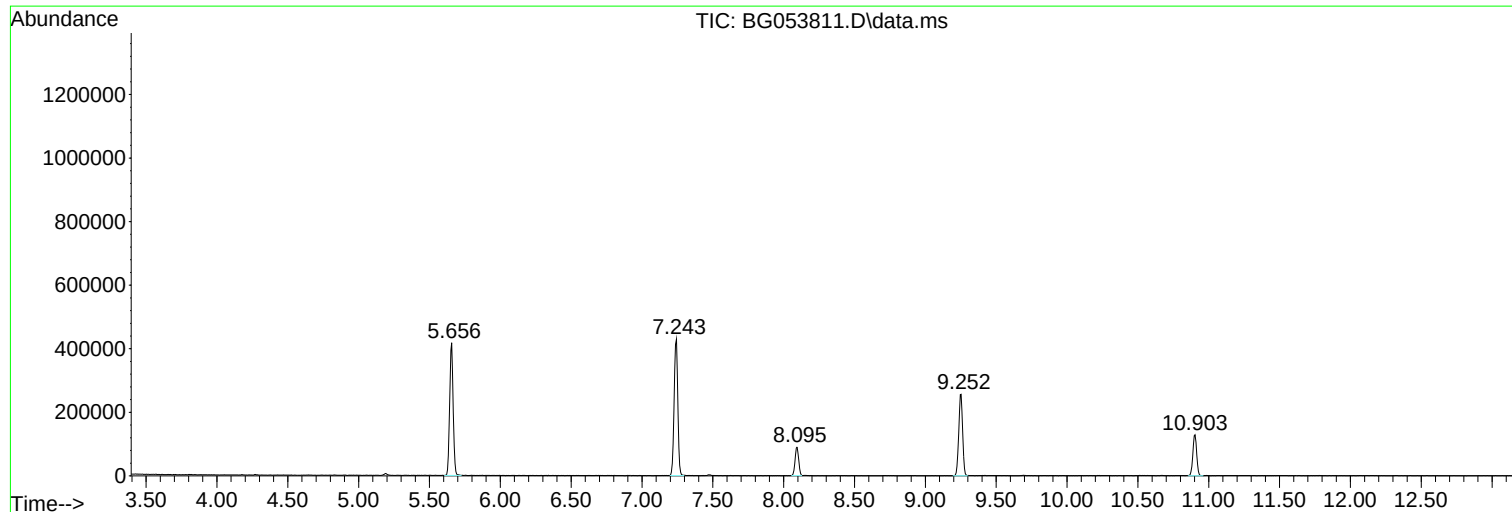
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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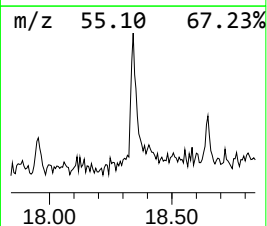
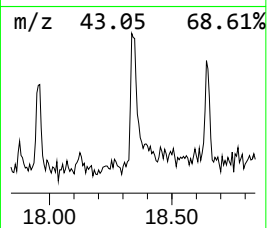
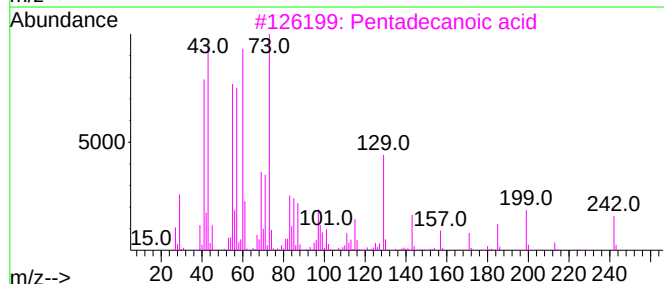
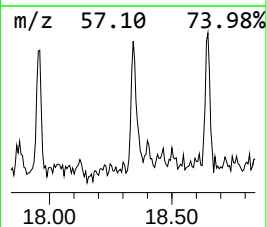
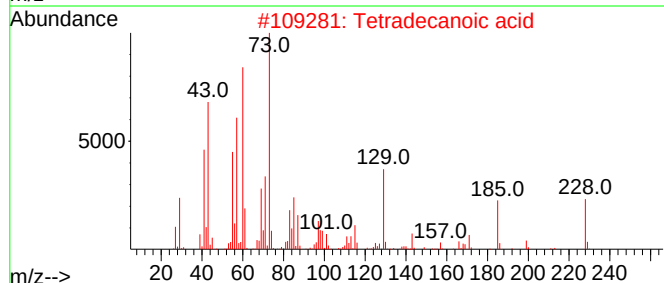
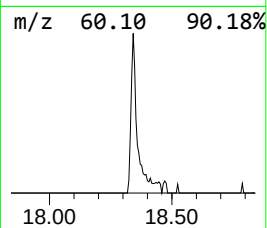
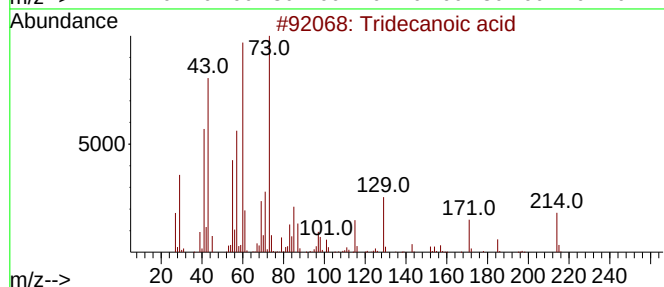
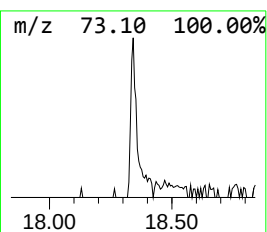
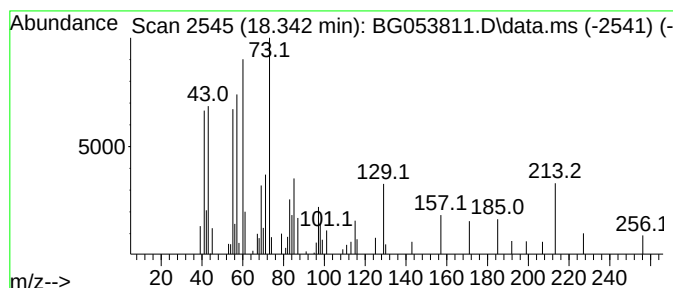
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Tridecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.342	2.34 ng	50448	Phenanthrene-d10	17.454

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanoic acid	214	C13H26O2	000638-53-9	58
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	58
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	58
4			Dodecanoic acid	200	C12H24O2	000143-07-7	58
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	55



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Misc :
ALS Vial : 8 Sample Multiplier: 1

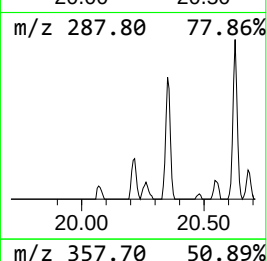
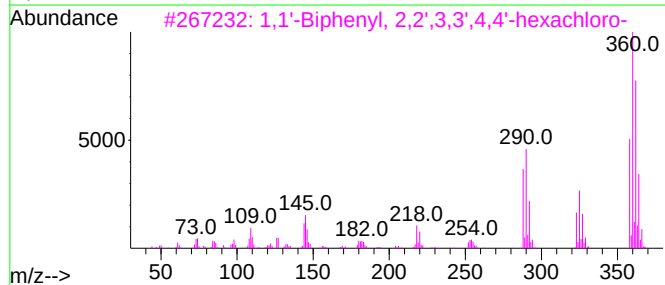
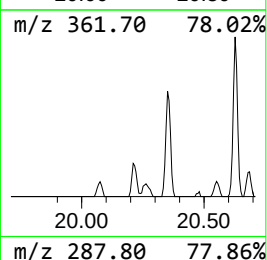
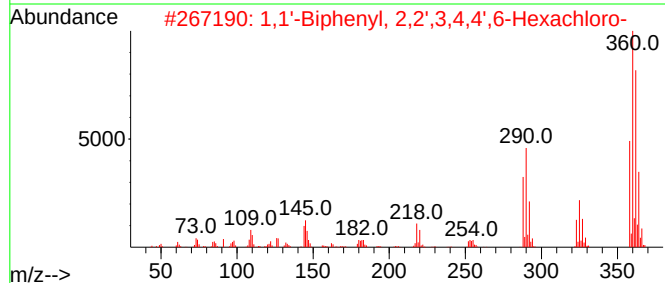
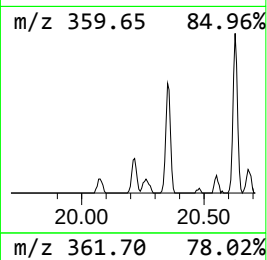
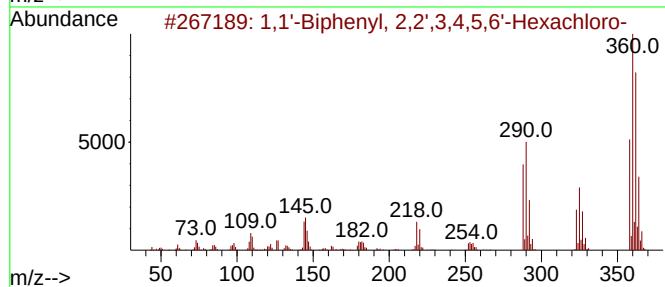
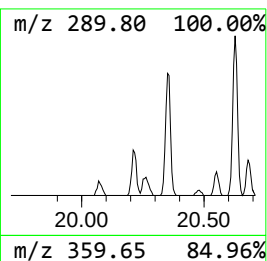
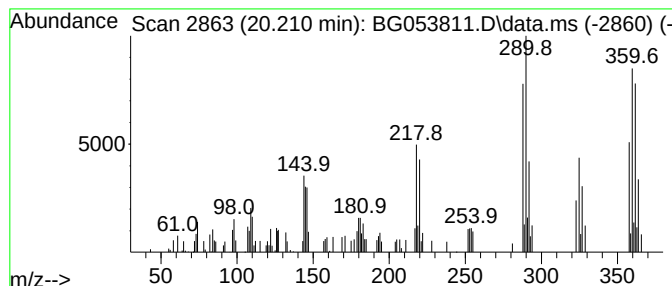
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 1,1'-Biphenyl, 2,2',3,4,5,6... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.210	2.20 ng	55100	Chrysene-d12	21.732	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,4,5,6'-Hex...	358	C12H4Cl6	068194-15-0	99
2	1,1'-Biphenyl, 2,2',3,4,4',6-Hex...	358	C12H4Cl6	056030-56-9	99
3	1,1'-Biphenyl, 2,2',3,3',4,4'-he...	358	C12H4Cl6	038380-07-3	99
4	1,1'-Biphenyl, 2,2',3,5,5',6-hex...	358	C12H4Cl6	052663-63-5	99
5	1,1'-Biphenyl, 2,2',3,3',5,5'-He...	358	C12H4Cl6	035694-04-3	99



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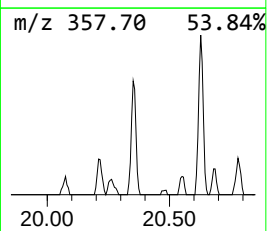
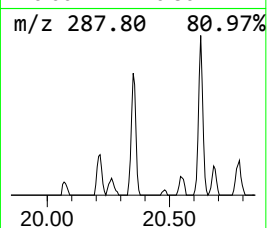
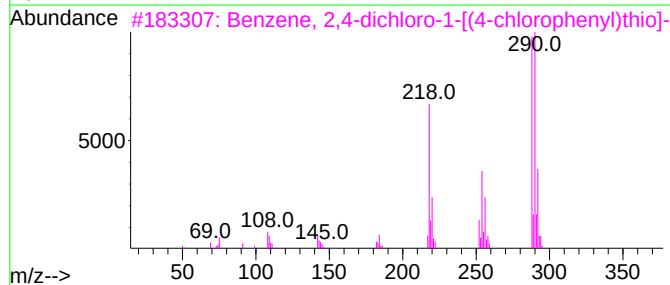
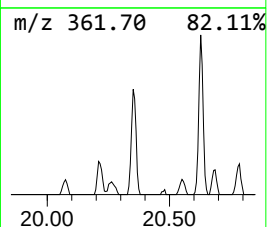
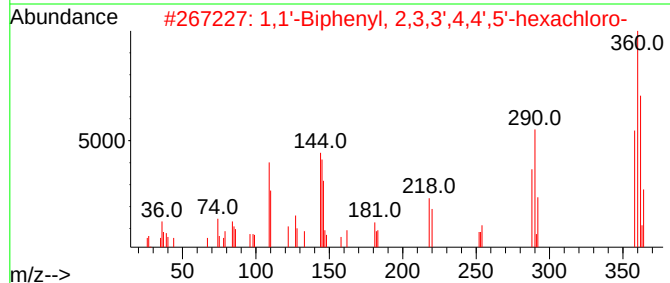
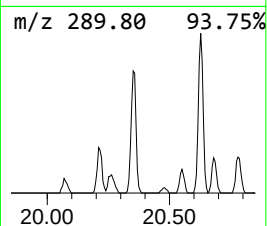
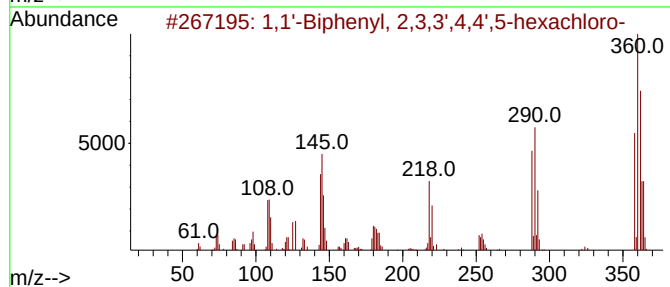
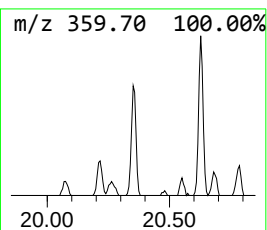
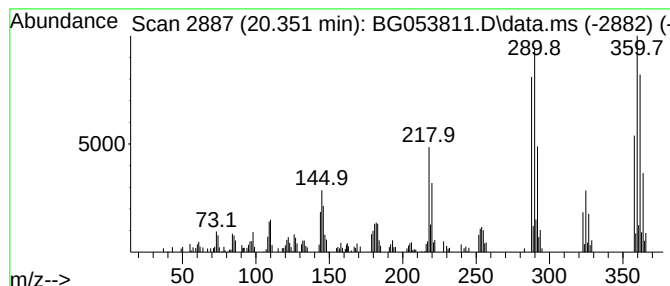
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.351	7.47 ng	187270	Chrysene-d12	21.732		
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl,	2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	91
2	1,1'-Biphenyl,	2,3,3',4,4',5'-he...	358	C12H4Cl6	069782-90-7	76
3	Benzene,	2,4-dichloro-1-[(4-chlo...	288	C12H7Cl3S	055759-88-1	56
4	1,3-Butadiyne,	1,4-di(4-bromophe...	358	C16H8Br2	000959-88-6	35
5	3-Bromo-5,7-dichloro-1-methylnap...		288	C11H7BrCl2	055058-84-9	30



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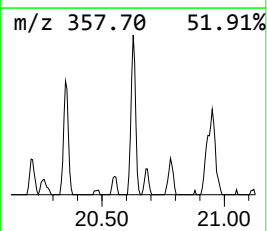
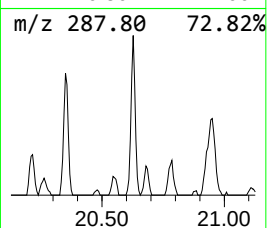
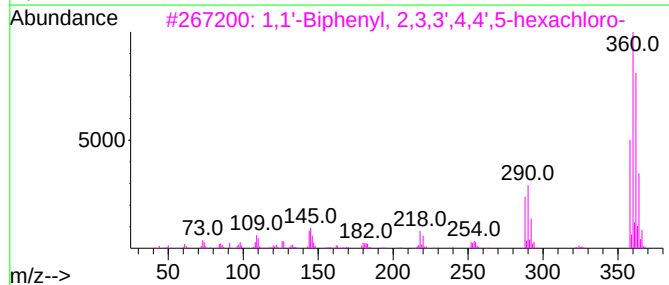
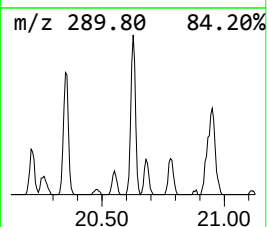
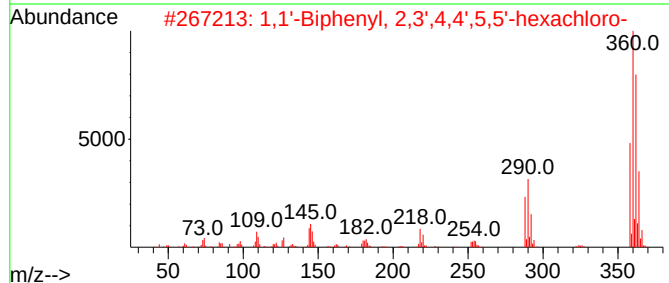
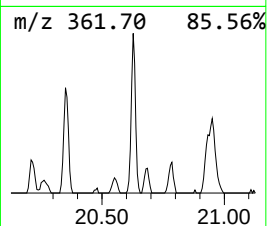
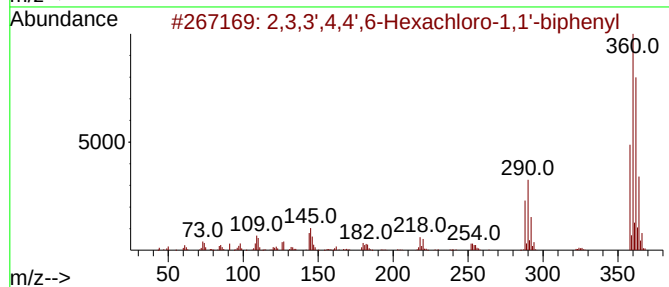
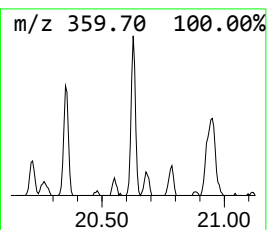
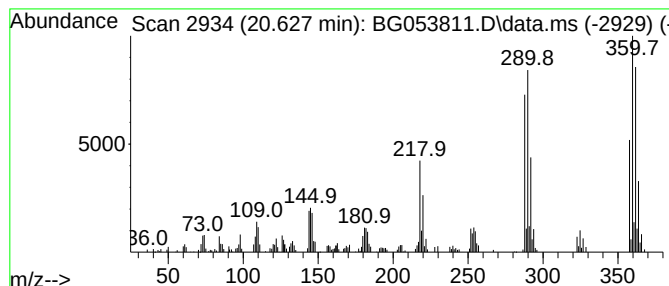
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.627	7.45 ng	186769	Chrysene-d12	21.732

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,3',4,4',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-42-7	99		
2	1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	99		
3	1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99		
4	2,3,3',4,5,5'-Hexachloro-1,1'-bi...	358	C12H4Cl6	039635-35-3	99		
5	2,3,3',4,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-43-8	99		



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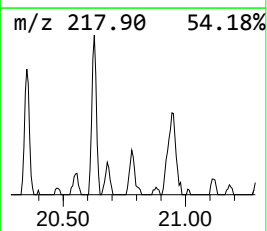
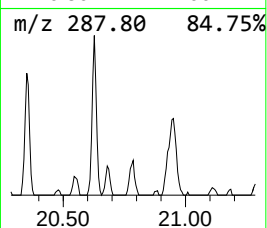
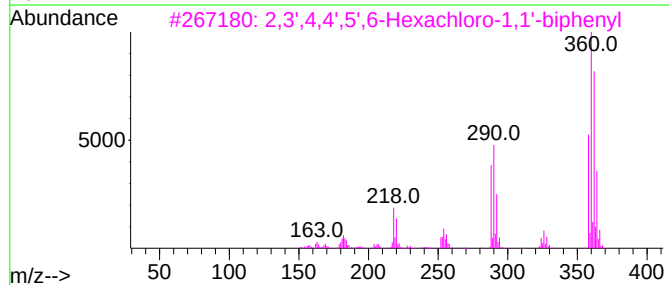
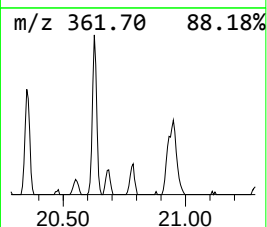
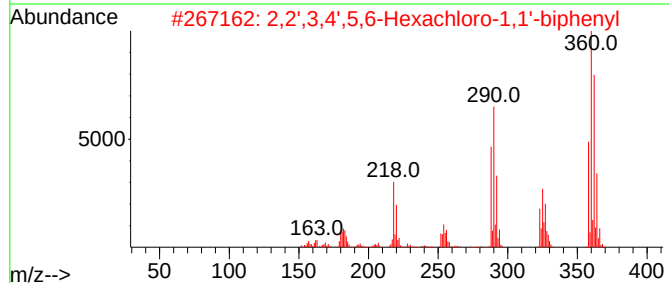
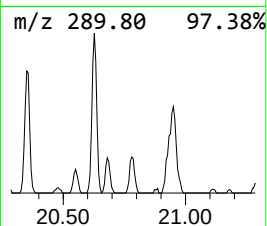
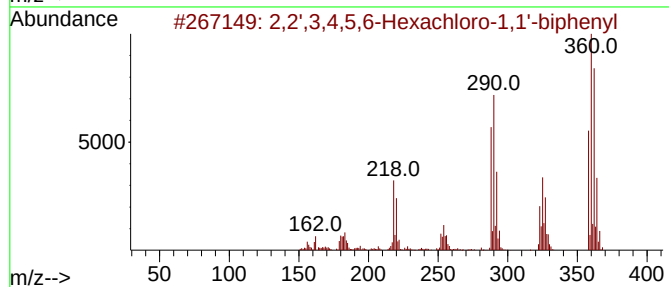
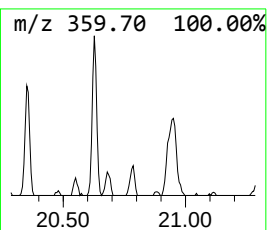
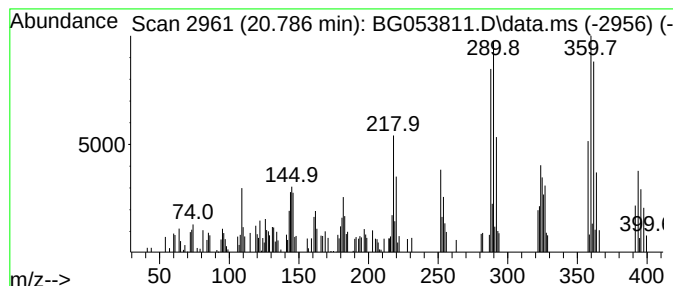
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.786	3.83 ng	96063	Chrysene-d12	21.732	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4,5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-61-4	94
2	2,2',3,4',5,6-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-13-8	93
3	2,3',4,4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	93
4	2,3,3',4',5,5'-Hexachloro-1,1'-b...	358	C12H4Cl6	039635-34-2	93
5	1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060322\
Data File : BG053811.D
Acq On : 3 Jun 2022 13:40
Operator : CG/JU
Sample : N2984-17
Misc :
ALS Vial : 8 Sample Multiplier: 1

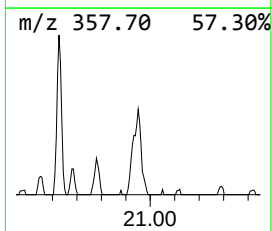
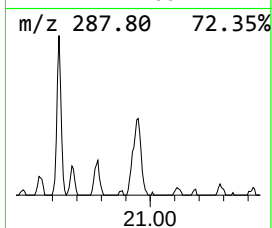
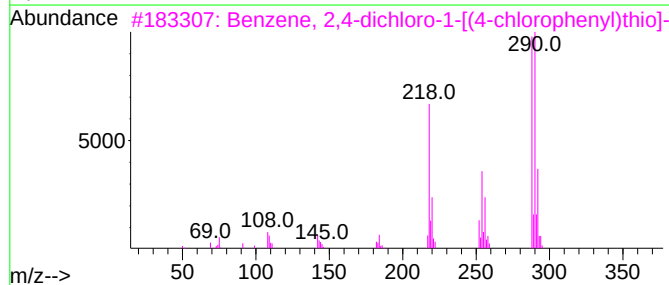
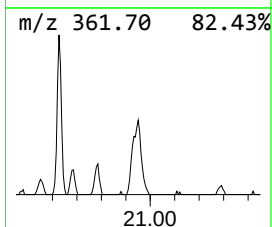
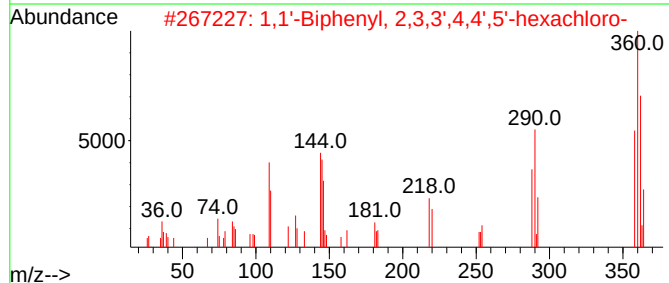
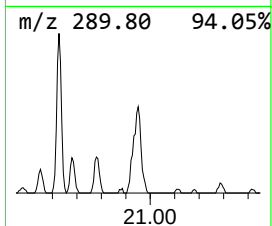
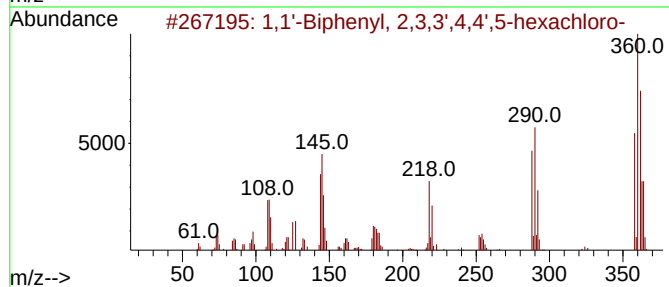
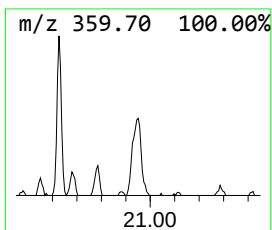
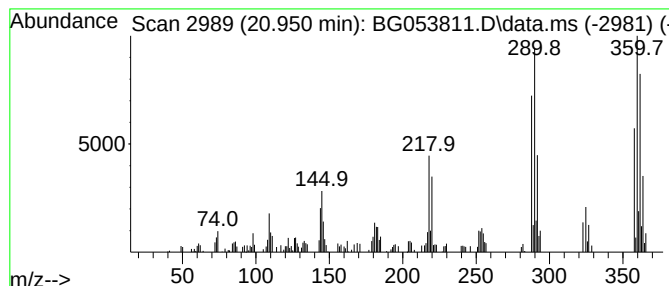
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.950	7.76 ng	194531	Chrysene-d12		21.732	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3',4,4',5-hex...		358	C12H4Cl6	038380-08-4	91
2	1,1'-Biphenyl, 2,3,3',4,4',5'-he...		358	C12H4Cl6	069782-90-7	87
3	Benzene, 2,4-dichloro-1-[(4-chlo...		288	C12H7Cl3S	055759-88-1	81
4	4,4'3'-Trichlorodiphenylsulfide		288	C12H7Cl3S	1000283-58-7	81
5	Penclomedine		323	C8H6Cl5NO2	108030-77-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060322\
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Sample : N2984-17
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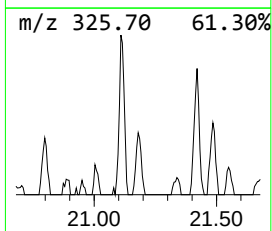
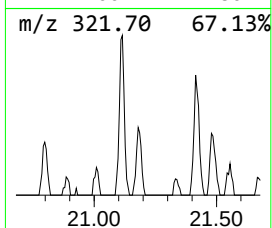
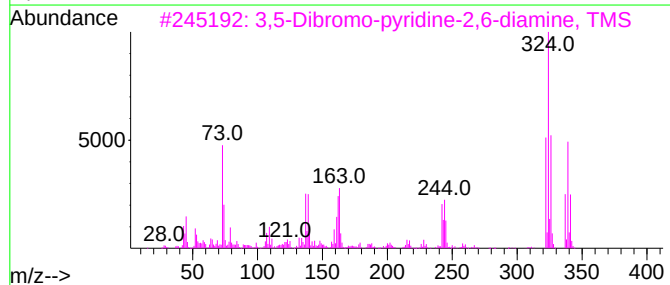
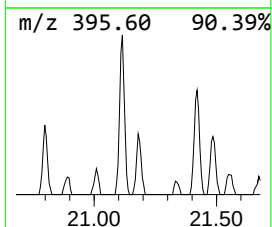
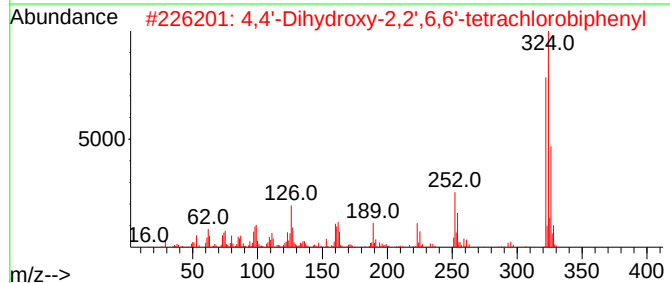
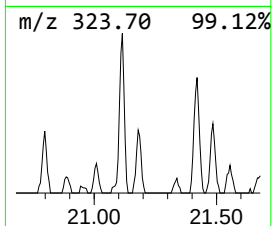
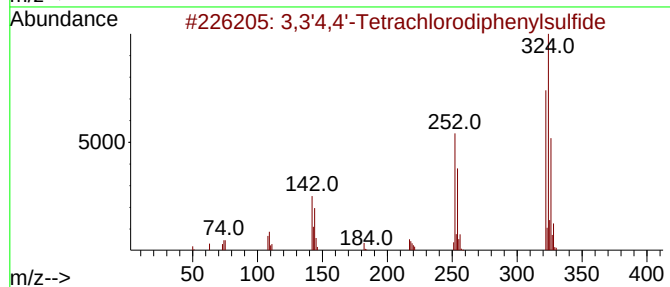
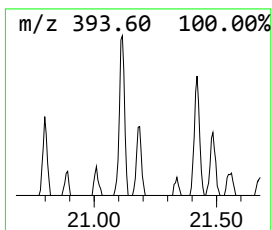
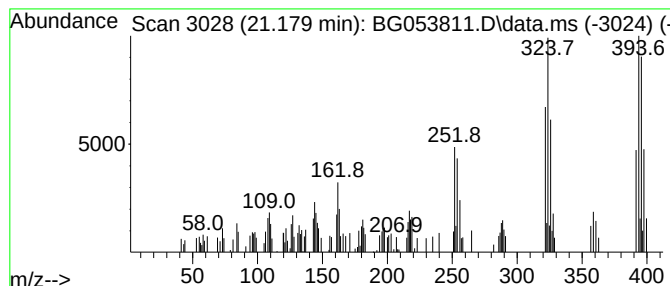
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 3,3',4,4'-Tetrachlorodipheny... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.180	2.19 ng	54860	Chrysene-d12	21.732	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3',4,4'-Tetrachlorodiphenylsulfide	322	C12H6Cl4S	076745-12-5	94
2	4,4'-Dihydroxy-2,2',6,6'-tetrach...	322	C12H6Cl4O2	1000447-97-4	38
3	3,5-Dibromo-pyridine-2,6-diamine...	337	C8H13Br2N3Si	1000511-29-8	14
4	2,2',3,4,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-57-2	11
5	2,2',3,5,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-55-0	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060322\
Data File : BG053811.D
Acq On : 3 Jun 2022 13:40
Operator : CG/JU
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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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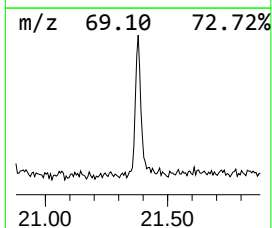
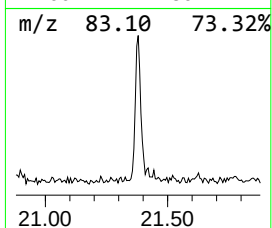
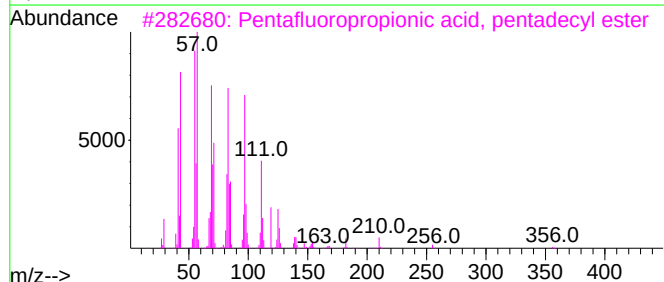
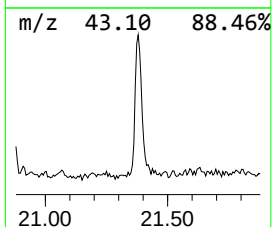
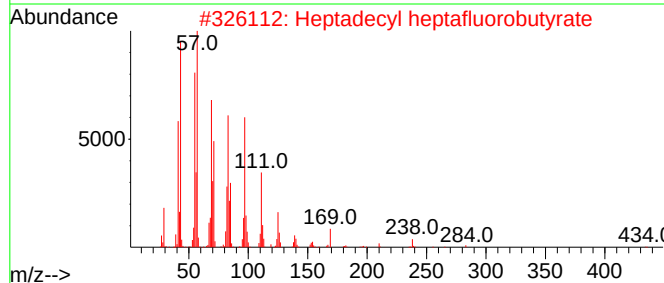
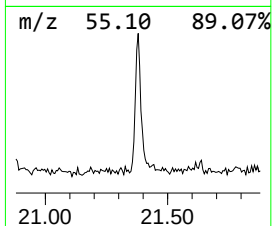
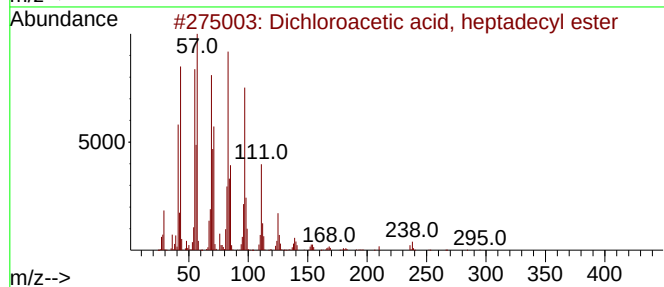
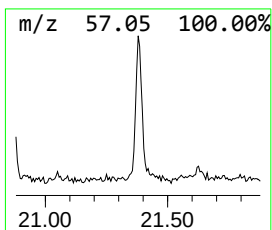
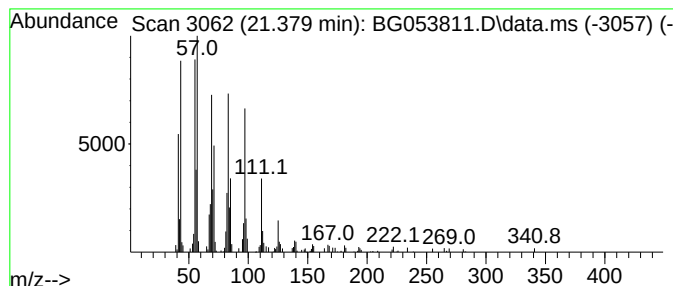
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Dichloroacetic acid, heptad... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.379	5.89 ng	147705	Chrysene-d12	21.732

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	95
2			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
4			Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
5			17-Pentatriacontene	491	C35H70	006971-40-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060322\
Data File : BG053811.D
Acq On : 3 Jun 2022 13:40
Operator : CG/JU
Sample : N2984-17
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ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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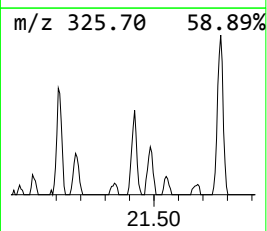
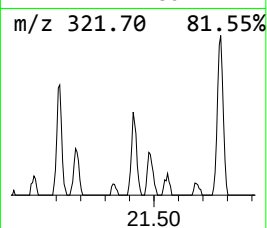
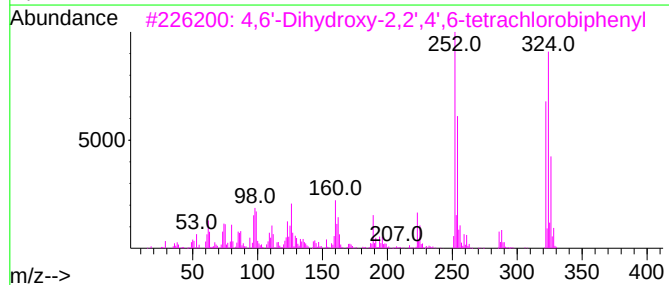
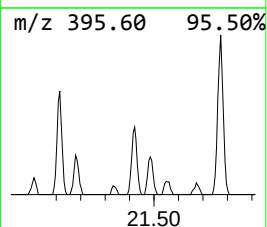
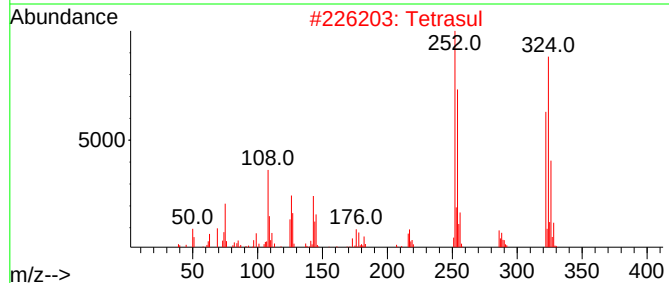
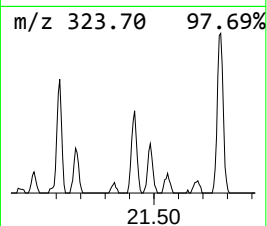
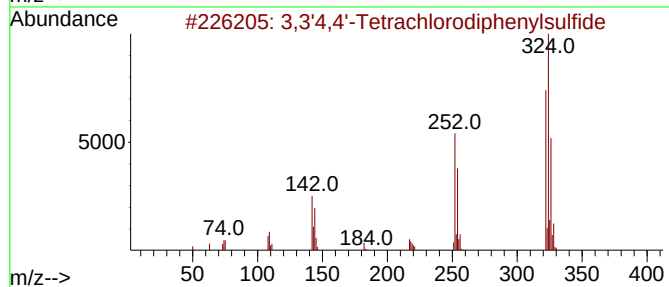
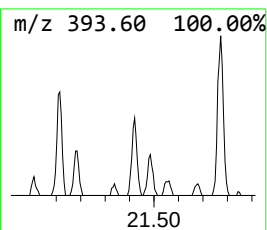
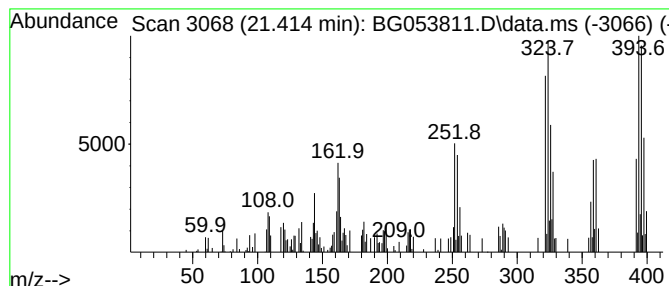
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown21.415 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.415	3.68 ng	92311	Chrysene-d12	21.732

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3',4,4'-Tetrachlorodiphenylsulfide	322	C12H6Cl4S	076745-12-5	52		
2	Tetrasul	322	C12H6Cl4S	002227-13-6	38		
3	4,6'-Dihydroxy-2,2',4',6-tetrach...	322	C12H6Cl4O2	1000447-97-5	30		
4	4,4'-Dihydroxy-2,2',6,6'-tetrach...	322	C12H6Cl4O2	1000447-97-4	25		
5	5,5'-Dibromo-2,2'-bithiophene	322	C8H4Br2S2	004805-22-5	14		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060322\
Data File : BG053811.D
Acq On : 3 Jun 2022 13:40
Operator : CG/JU
Sample : N2984-17
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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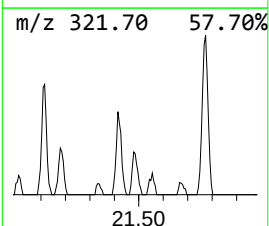
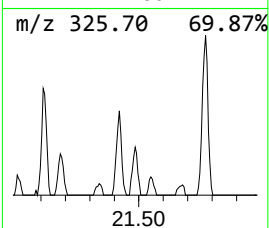
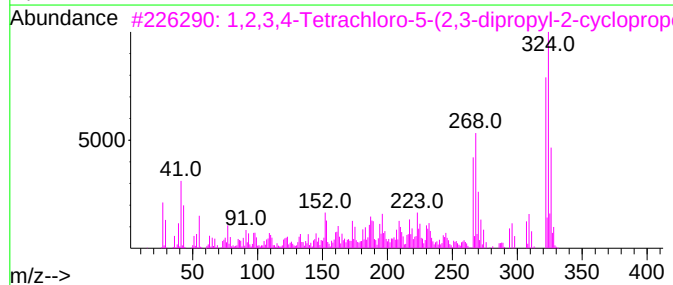
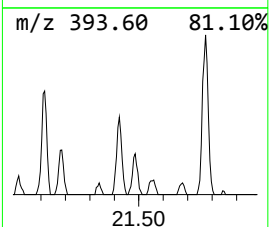
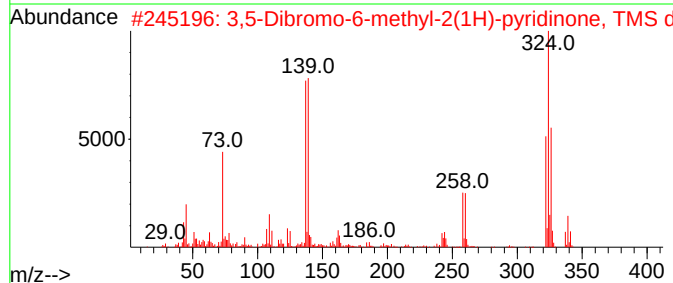
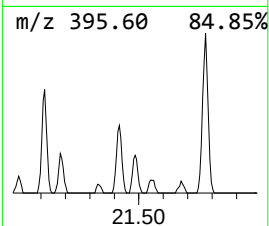
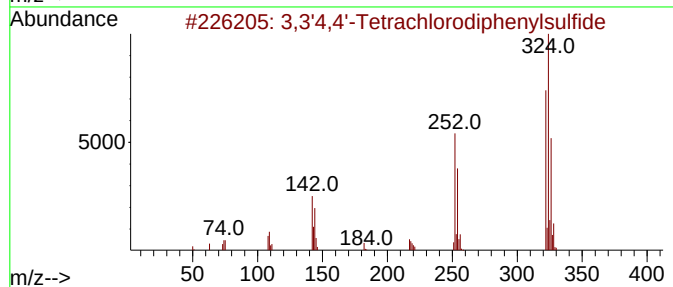
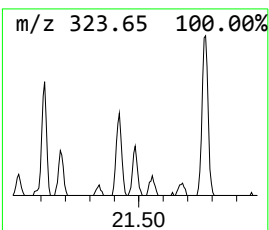
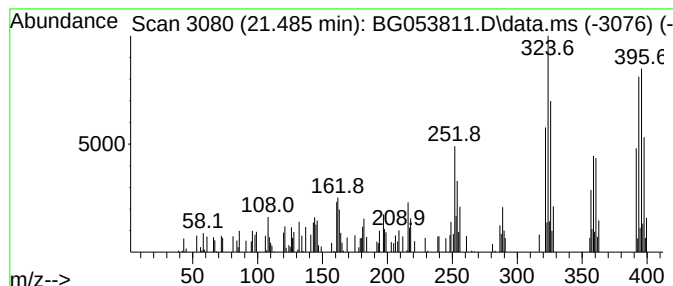
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown21.485 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.485	2.52 ng	63083	Chrysene-d12	21.732

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3',4,4'-Tetrachlorodiphenylsulfide	322	C12H6Cl4S	076745-12-5	60		
2	3,5-Dibromo-6-methyl-2(1H)-pyrid...	337	C9H13Br2NOSi	1000474-24-8	14		
3	1,2,3,4-Tetrachloro-5-(2,3-dipro...	322	C14H14Cl4	005680-25-1	11		
4	4,4'-Dihydroxy-2,2',6,6'-tetrach...	322	C12H6Cl4O2	1000447-97-4	11		
5	1,3,7-Trichloro-9-phenanthroic acid	324	C15H7Cl3O2	056988-57-9	11		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060322\
Data File : BG053811.D
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Operator : CG/JU
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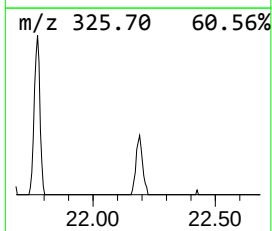
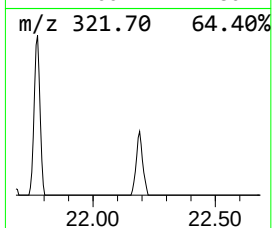
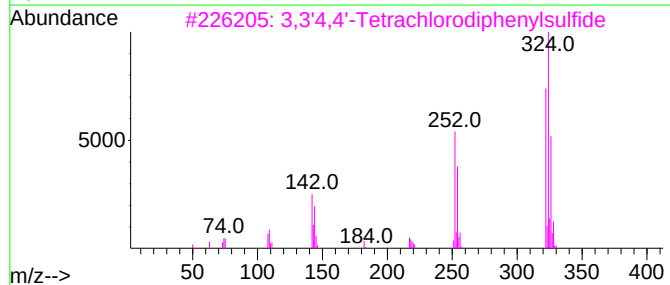
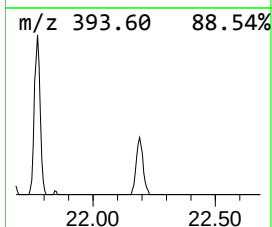
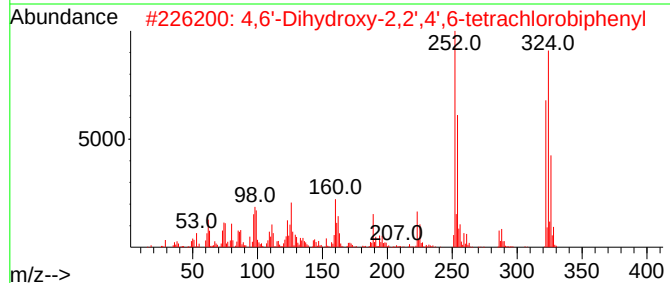
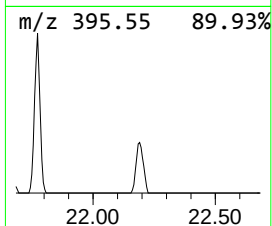
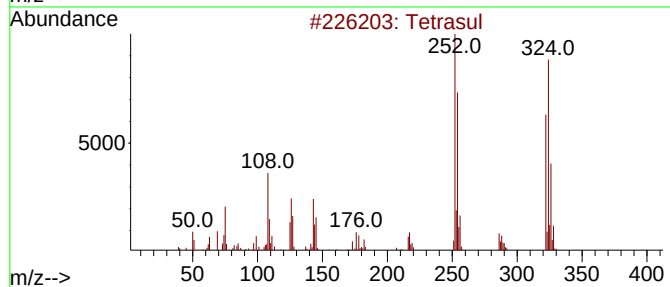
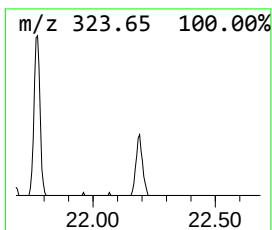
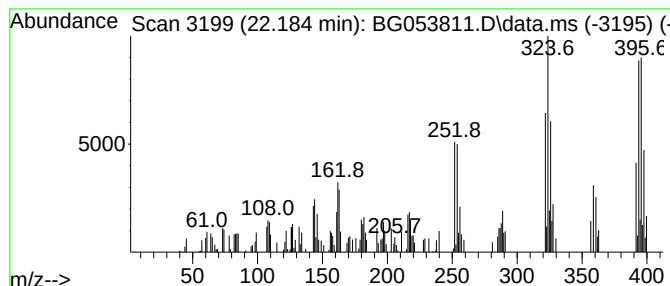
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown22.184 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.184	4.10 ng	102702	Chrysene-d12	21.732

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrasul	322	C12H6Cl4S	002227-13-6	41
2			4,6'-Dihydroxy-2,2',4',6-tetrach...	322	C12H6Cl4O2	1000447-97-5	38
3			3,3',4,4'-Tetrachlorodiphenylsulfide	322	C12H6Cl4S	076745-12-5	38
4			4,4'-Dihydroxy-2,2',6,6'-tetrach...	322	C12H6Cl4O2	1000447-97-4	25
5			3,5-Dibromo-pyridine-2,6-diamine...	337	C8H13Br2N3Si	1000511-29-8	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060322\
Data File : BG053811.D
Acq On : 3 Jun 2022 13:40
Operator : CG/JU
Sample : N2984-17
Misc :
ALS Vial : 8 Sample Multiplier: 1

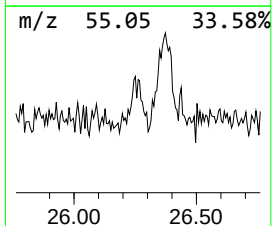
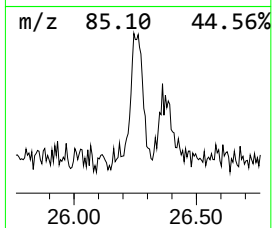
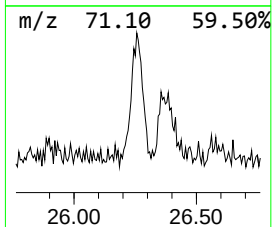
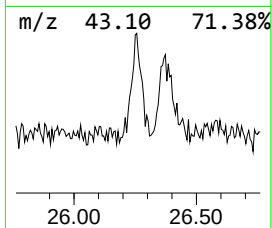
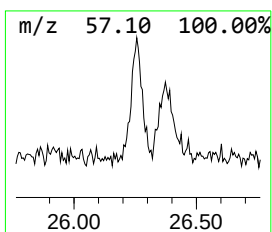
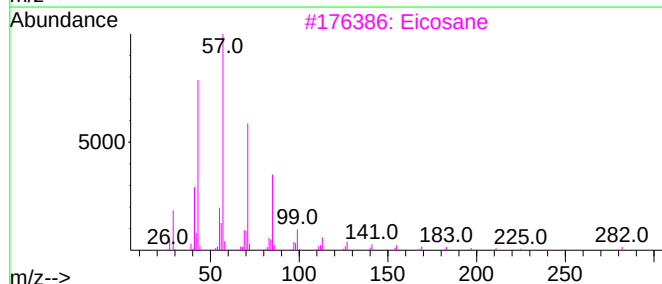
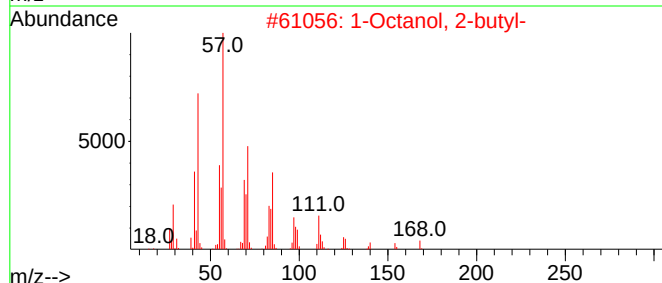
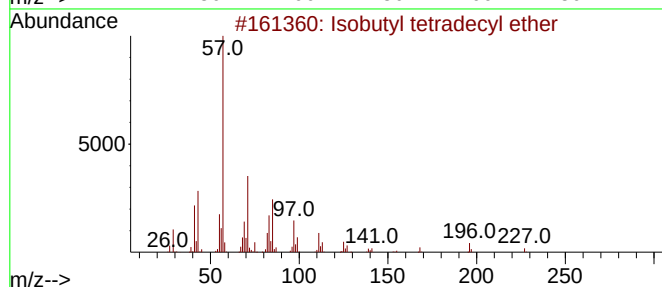
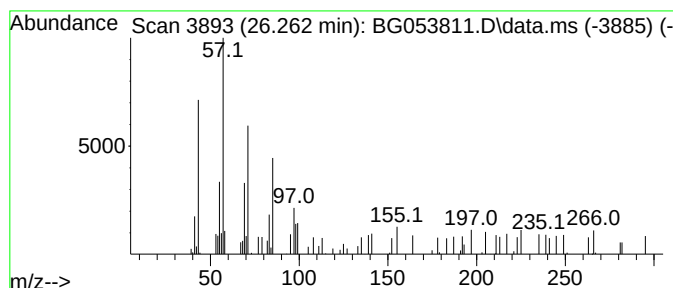
Instrument :
BNA_G
ClientSampleId :
P021-SS001-0612-01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Isobutyl tetradecyl ether Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.262	2.10 ng	48399	Perylene-d12	25.004	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isobutyl tetradecyl ether	270	C18H38O	1000406-32-7	50
2	1-Octanol, 2-butyl-	186	C12H26O	003913-02-8	50
3	Eicosane	282	C20H42	000112-95-8	47
4	Oct-3-enoylamide, N-undecyl-	295	C19H37NO	1010420-41-6	47
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060322\
Data File : BG053811.D
Acq On : 3 Jun 2022 13:40
Operator : CG/JU
Sample : N2984-17
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P021-SS001-0612-01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Tridecanoic acid	18.342	2.3	ng	50448	4	17.454	430517	20.0
1,1'-Biphenyl, ...	20.210	2.2	ng	55100	5	21.732	501120	20.0
1,1'-Biphenyl, ...	20.351	7.5	ng	187270	5	21.732	501120	20.0
2,3,3',4,4',6-H...	20.627	7.5	ng	186769	5	21.732	501120	20.0
2,2',3,4,5,6-He...	20.786	3.8	ng	96063	5	21.732	501120	20.0
1,1'-Biphenyl, ...	20.950	7.8	ng	194531	5	21.732	501120	20.0
1,1'-Biphenyl, ...	21.115	5.2	ng	130579	5	21.732	501120	20.0
3,3'4,4'-Tetrac...	21.180	2.2	ng	54860	5	21.732	501120	20.0
Dichloroacetic ...	21.379	5.9	ng	147705	5	21.732	501120	20.0
unknown21.415	21.415	3.7	ng	92311	5	21.732	501120	20.0
unknown21.485	21.485	2.5	ng	63083	5	21.732	501120	20.0
unknown22.184	22.184	4.1	ng	102702	5	21.732	501120	20.0
Isobutyl tetrad...	26.262	2.1	ng	48399	6	25.004	460676	20.0