

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
 Data File : BM037442.D
 Acq On : 09 Nov 2022 18:16
 Operator : CG/JU
 Sample : N5436-05 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 250FERRYBLVD-TFS-05

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_M\Methods\8270-BM102822.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM037442.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.934	308	314	320	rBV	140853	187134	1.55%	0.283%
2	5.399	386	393	398	rBV	341427	505678	4.19%	0.766%
3	7.005	661	666	682	rVB	302691	528486	4.38%	0.800%
4	7.822	799	805	813	rVB	1102084	1653874	13.70%	2.504%
5	8.604	933	938	946	rBV	76845	143452	1.19%	0.217%
6	8.993	998	1004	1008	rBV	191131	327142	2.71%	0.495%
7	9.810	1134	1143	1161	rBV	141945	353498	2.93%	0.535%
8	10.051	1174	1184	1192	rBV5	48780	154066	1.28%	0.233%
9	10.622	1274	1281	1286	rBV	1384139	2261282	18.73%	3.424%
10	10.669	1286	1289	1296	rVB	190670	292732	2.43%	0.443%
11	13.086	1693	1700	1707	rBV	531073	773462	6.41%	1.171%
12	13.298	1728	1736	1742	rBV3	55492	127639	1.06%	0.193%
13	13.945	1837	1846	1854	rBV4	88406	180189	1.49%	0.273%
14	14.186	1881	1887	1897	rBV	83575	137935	1.14%	0.209%
15	14.463	1925	1934	1940	rBV	2212075	3157438	26.16%	4.781%
16	14.522	1940	1944	1952	rVB	131808	203005	1.68%	0.307%
17	14.857	1996	2001	2010	rBV	85803	150643	1.25%	0.228%
18	15.504	2106	2111	2122	rBV	132731	217965	1.81%	0.330%
19	15.710	2142	2146	2151	rBV4	98910	141533	1.17%	0.214%
20	15.951	2182	2187	2193	rBV	443955	625766	5.18%	0.948%
21	17.204	2394	2400	2404	rBV	2663201	3722535	30.84%	5.637%
22	17.245	2404	2407	2415	rVV	515251	744976	6.17%	1.128%
23	17.333	2418	2422	2426	rVV	180534	231729	1.92%	0.351%
24	17.804	2497	2502	2505	rBV4	78604	153585	1.27%	0.233%
25	17.851	2505	2510	2517	rVB	1978217	2404204	19.92%	3.641%
26	18.098	2547	2552	2567	rBV3	921084	1637666	13.57%	2.480%
27	18.274	2578	2582	2586	rBV2	184696	287274	2.38%	0.435%
28	18.386	2597	2601	2606	rBV2	220982	352912	2.92%	0.534%
29	18.604	2633	2638	2643	rBV2	625336	880391	7.29%	1.333%
30	18.980	2699	2702	2704	rBV2	199306	249864	2.07%	0.378%
31	19.015	2704	2708	2718	rVB	688685	1046579	8.67%	1.585%
32	19.263	2742	2750	2755	rBV	1446213	2515496	20.84%	3.809%
33	19.310	2755	2758	2765	rVB	423131	530079	4.39%	0.803%
34	19.374	2766	2769	2773	rBV	635964	893902	7.41%	1.354%
35	19.621	2803	2811	2821	rBV	9709079	12070476	100.00%	18.277%
36	19.827	2842	2846	2848	rBV	854174	996315	8.25%	1.509%

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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_M\Methods\8270-BM102822.M
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37	19.862	2848	2852	2859	rVB	5126582	6170079	51.12%	9.343%
38	20.268	2917	2921	2926	rBV2	358512	556708	4.61%	0.843%
39	20.551	2966	2969	2971	rBV2	257407	322816	2.67%	0.489%
40	20.586	2972	2975	2979	rVB	524393	579726	4.80%	0.878%
41	20.845	3016	3019	3024	rVB2	895737	984559	8.16%	1.491%
42	21.004	3043	3046	3049	rBV	446144	566009	4.69%	0.857%
43	21.092	3058	3061	3064	rBV	567273	704763	5.84%	1.067%
44	21.380	3105	3110	3121	rVB	3559011	6493290	53.79%	9.832%
45	21.774	3174	3177	3185	rVB	1158945	1481874	12.28%	2.244%
46	21.845	3186	3189	3192	rBV	1137394	1402570	11.62%	2.124%
47	22.433	3287	3289	3294	rVB2	593841	777785	6.44%	1.178%
48	22.903	3366	3369	3375	rVB2	808495	1188888	9.85%	1.800%
49	23.721	3503	3508	3515	rVB2	2239374	3970400	32.89%	6.012%

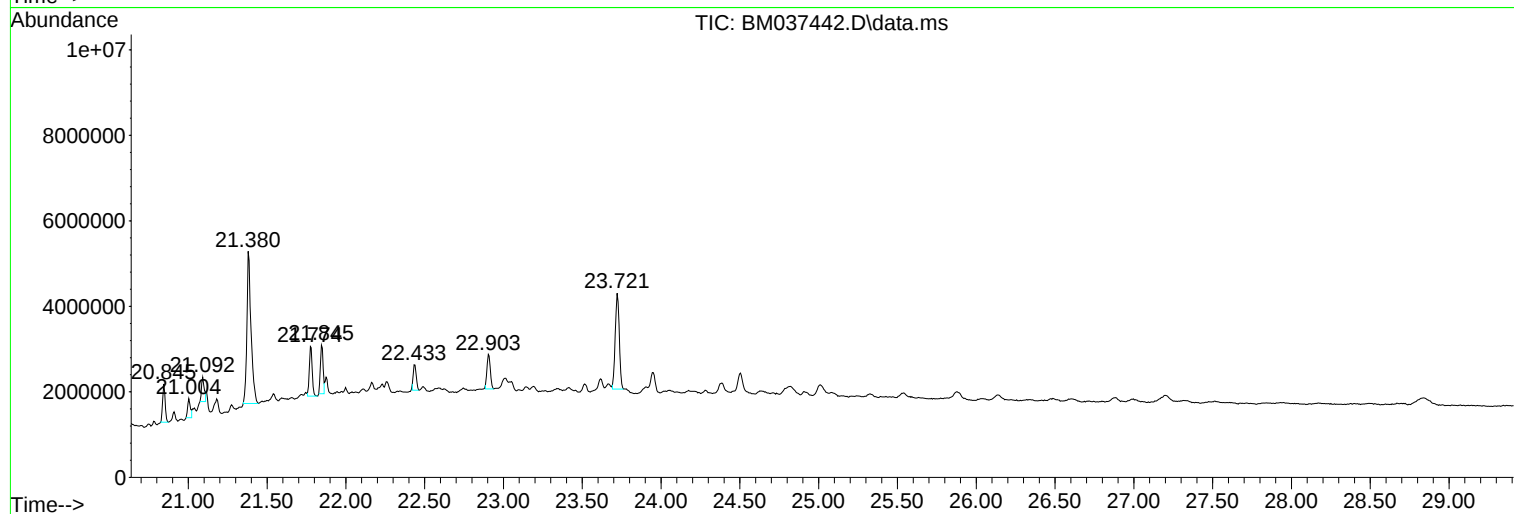
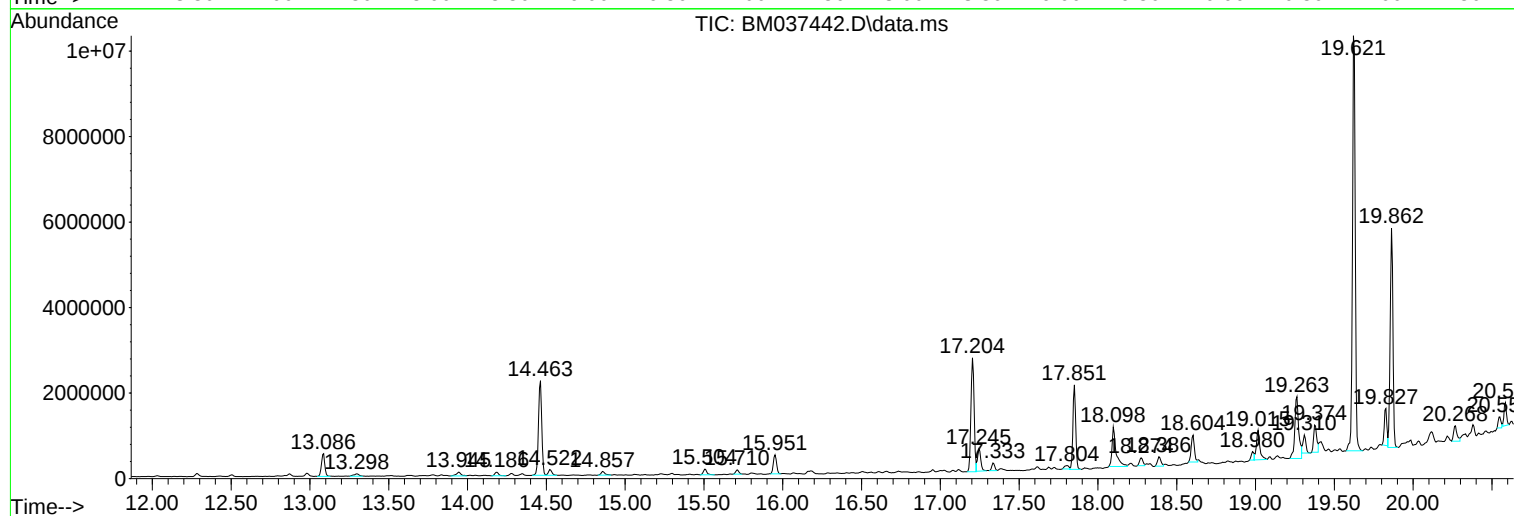
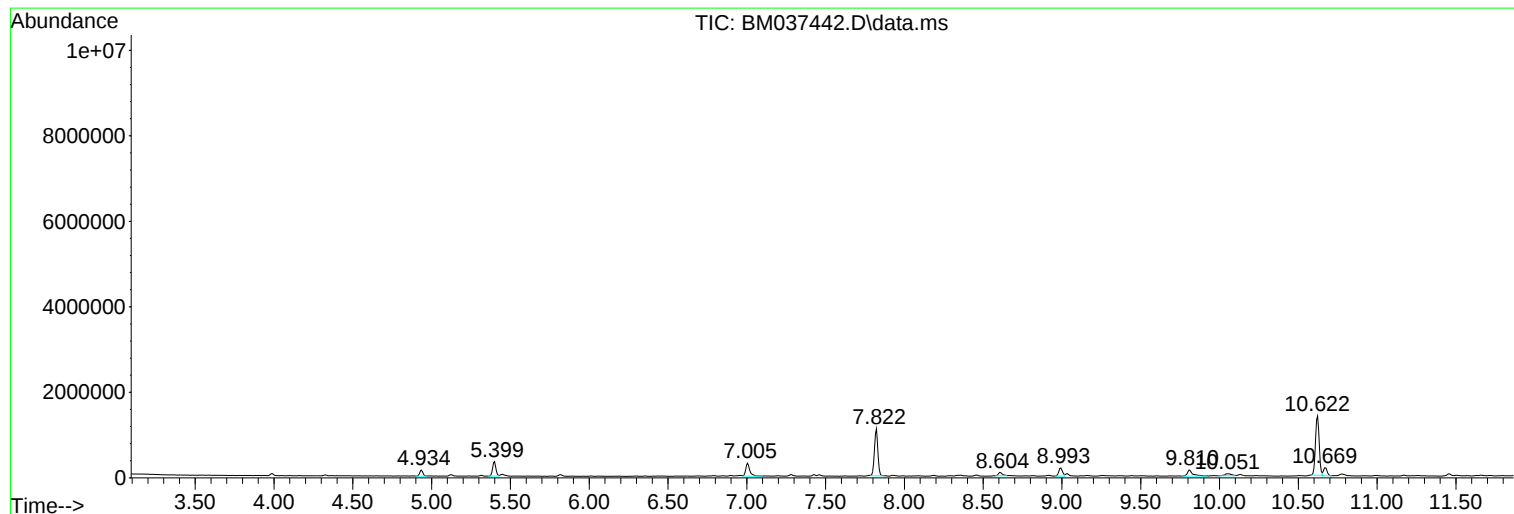
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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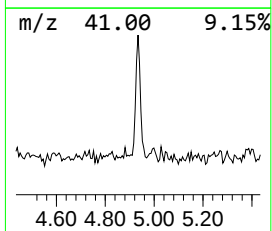
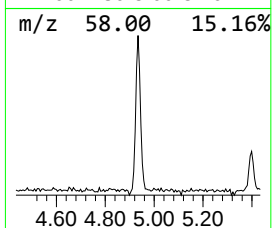
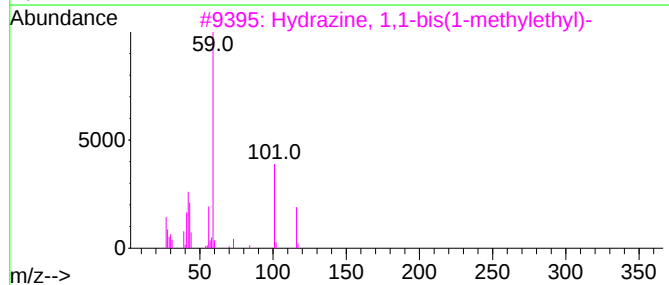
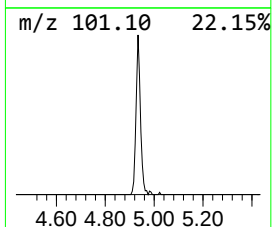
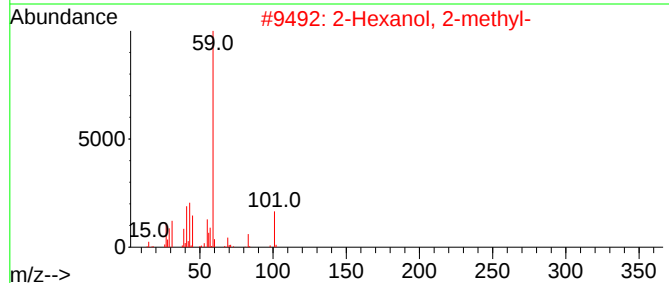
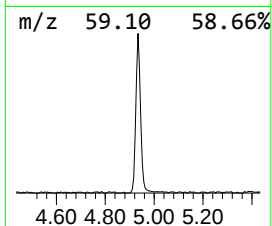
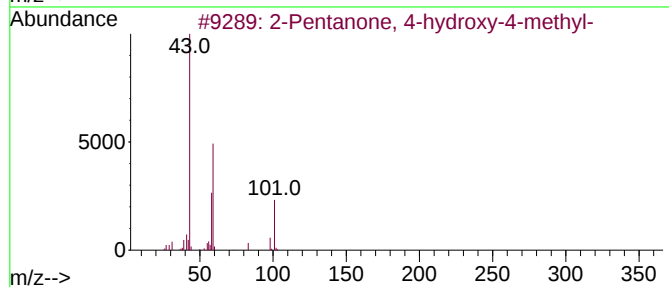
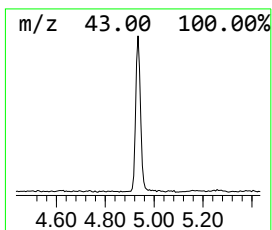
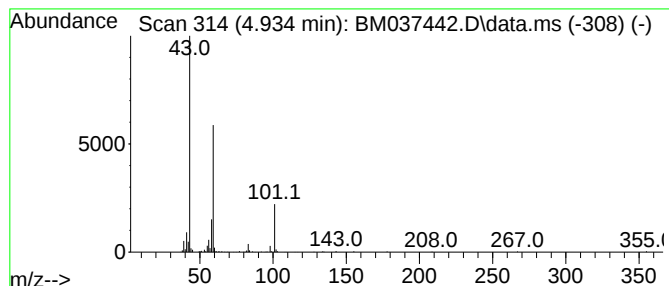
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.934	2.26 ng	187134	1,4-Dichlorobenzene-d4	7.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
3			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	28
4			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



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Instrument :
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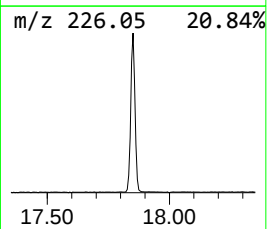
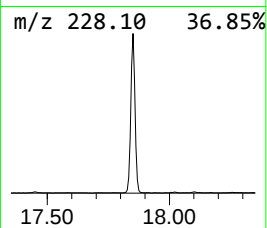
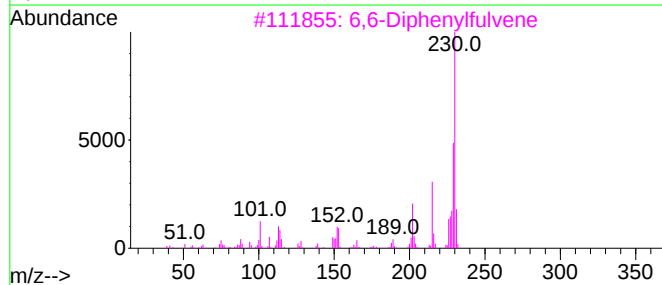
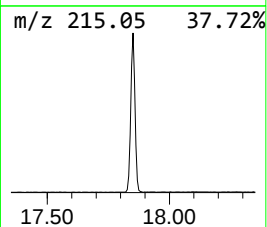
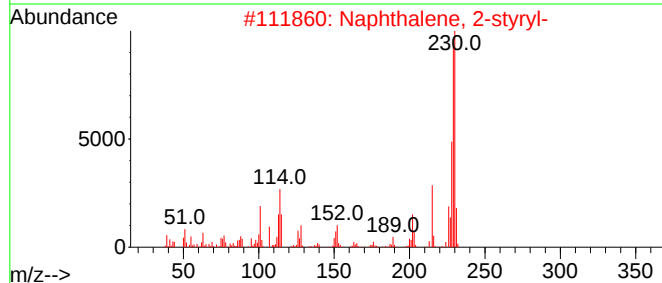
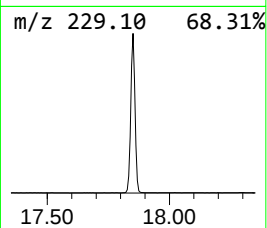
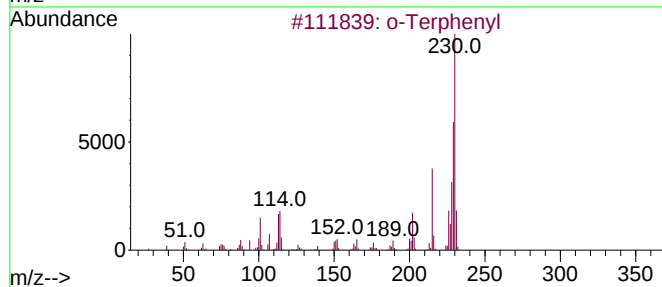
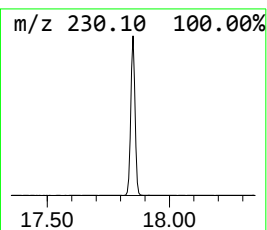
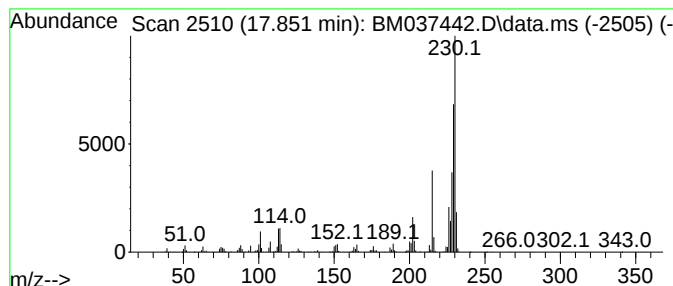
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 o-Terphenyl Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.851	12.92 ng	2404200	Phenanthrene-d10	17.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Terphenyl	230	C18H14	000084-15-1	98
2			Naphthalene, 2-styryl-	230	C18H14	002039-70-5	93
3			6,6-Diphenylfulvene	230	C18H14	002175-90-8	92
4			Pyrene, 1,3-dimethyl-	230	C18H14	064401-21-4	86
5			5,6-Dihydrochrysene	230	C18H14	002091-92-1	72



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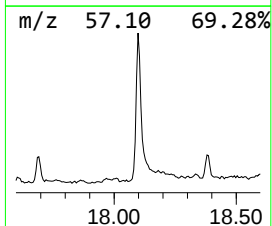
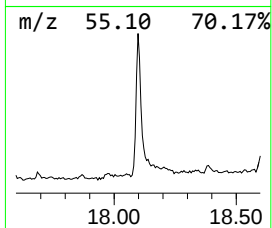
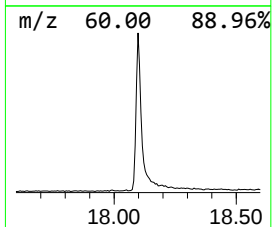
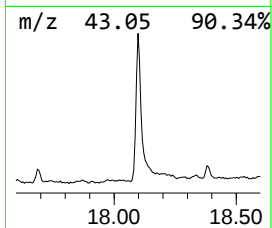
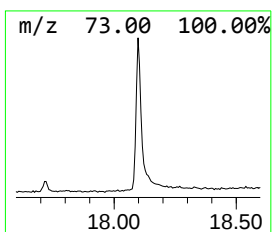
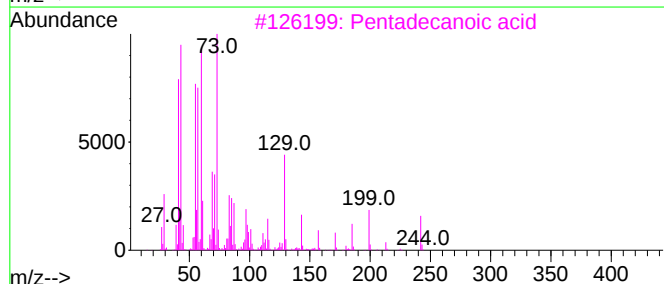
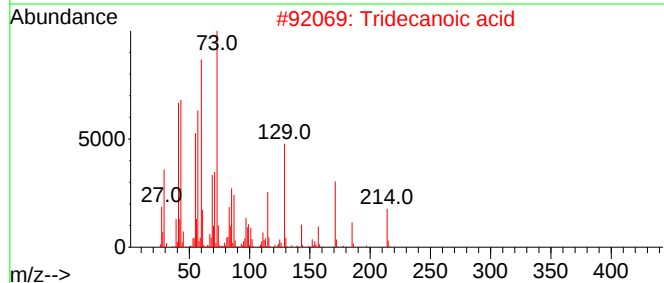
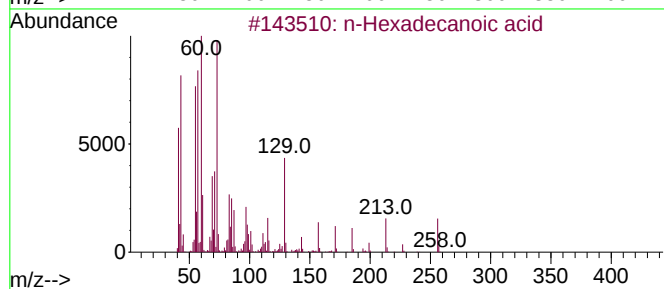
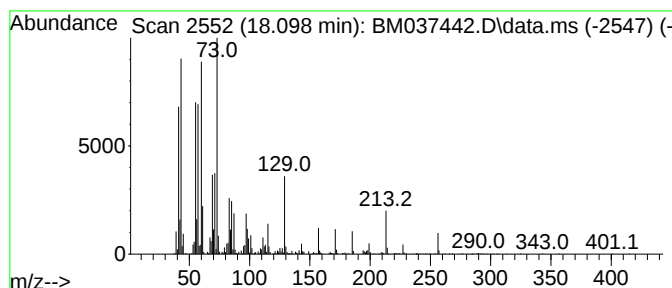
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.098	8.80 ng	1637670	Phenanthrene-d10	17.204	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	94
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5	Octadecanoic acid	284	C18H36O2	000057-11-4	60



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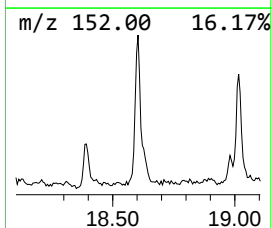
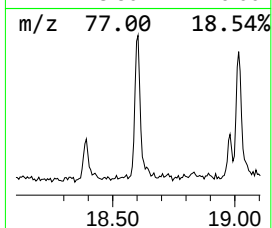
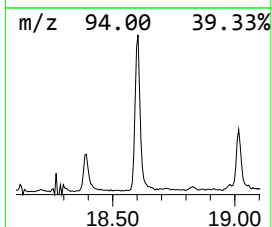
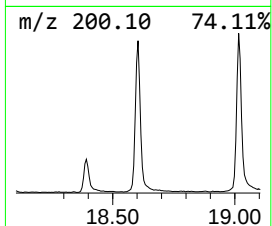
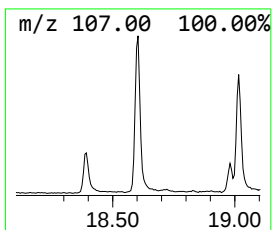
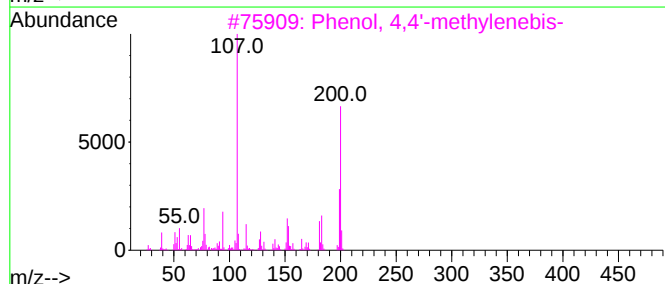
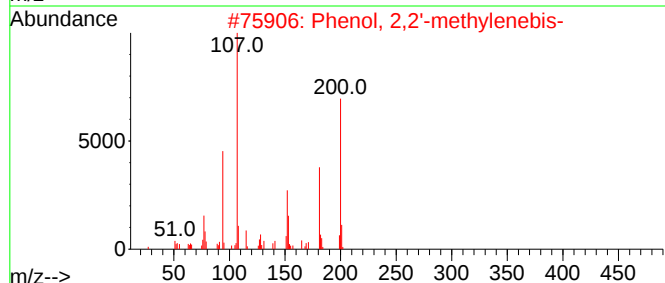
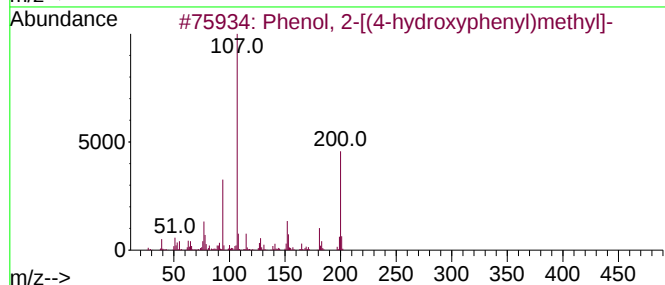
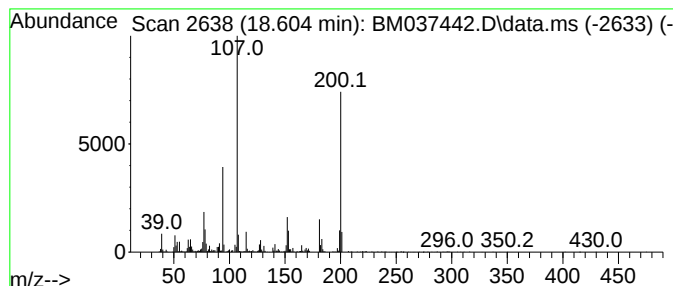
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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Peak Number 4 Phenol, 2-[(4-hydroxyphenyl)... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.604	4.73 ng	880391	Phenanthrene-d10	17.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	96
2			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	94
3			Phenol, 4,4'-methylenebis-	200	C13H12O2	000620-92-8	81
4			2,4'-Dihydroxydiphenylmethane, d...	284	C17H16O4	959033-36-4	72
5			2,2'-Methylenediphenol, acetate	242	C15H14O3	1000462-45-6	52



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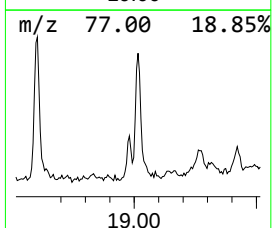
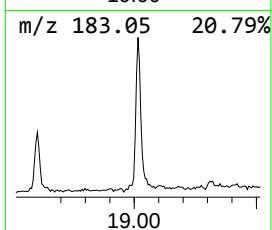
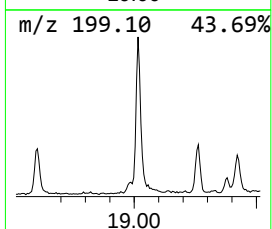
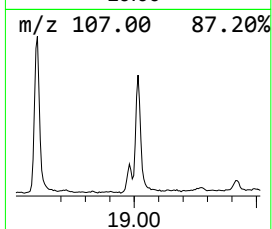
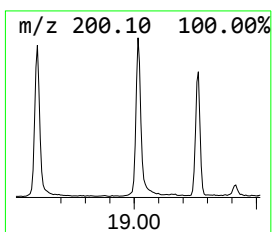
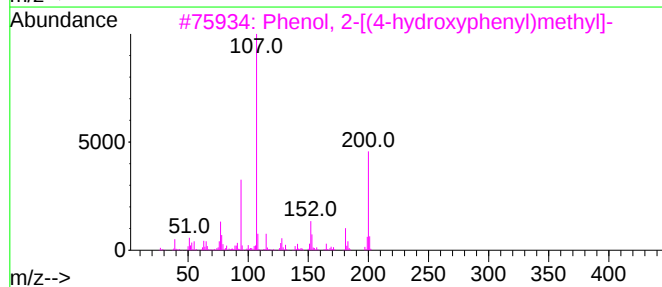
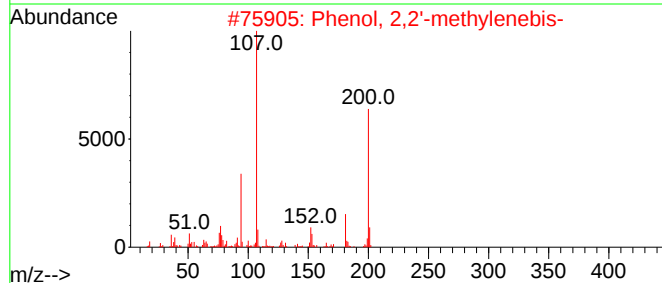
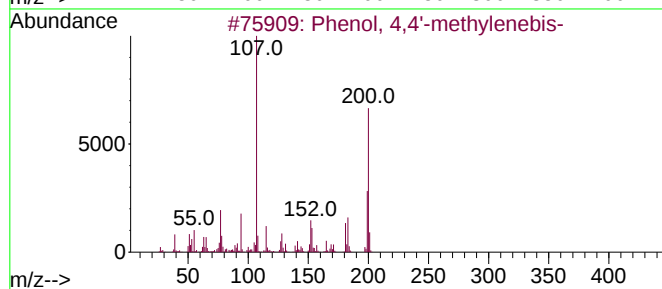
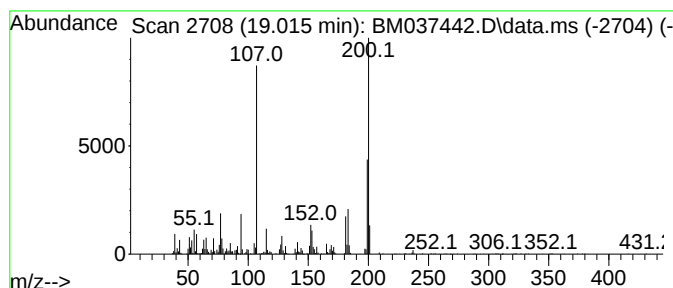
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BNA_M
ClientSampleId :
250FERRYBLVD-TFS-05

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Phenol, 4,4'-methylenebis- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.015	5.62 ng	1046580	Phenanthrene-d10	17.204	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-methylenebis-	200	C13H12O2	000620-92-8	98
2	Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	81
3	Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	81
4	Bisphenol F, 2-methylpropionate	270	C17H18O3	1000462-19-4	72
5	2,4'-Dihydroxydiphenylmethane, d...	284	C17H16O4	959033-36-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037442.D
Acq On : 09 Nov 2022 18:16
Operator : CG/JU
Sample : N5436-05 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-05

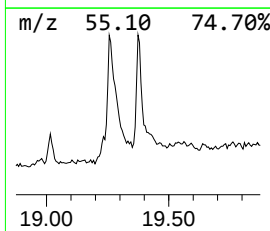
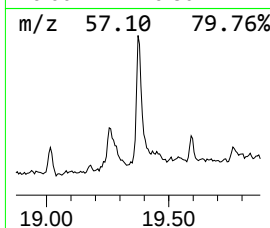
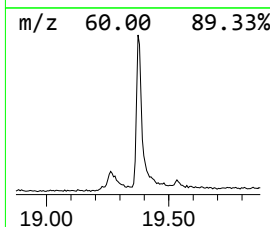
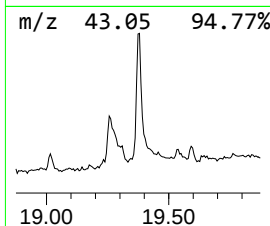
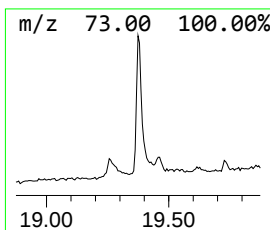
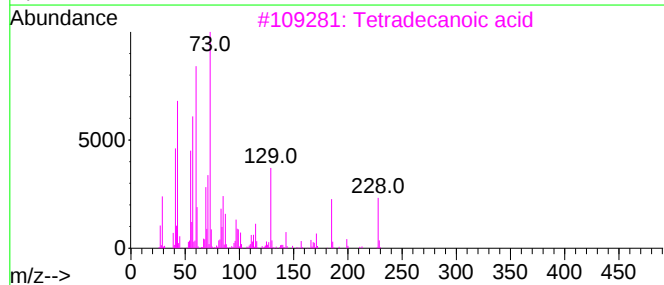
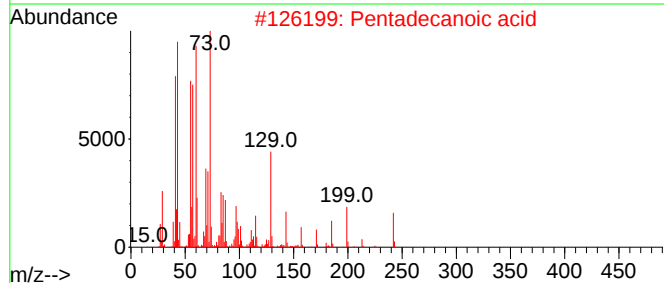
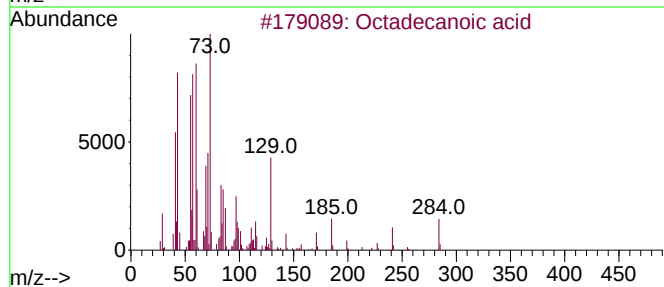
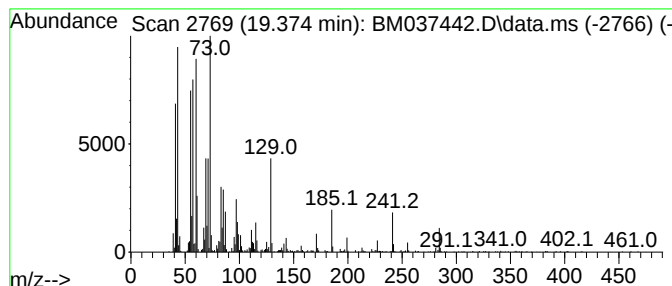
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.374	2.75 ng	893902	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4			n-Decanoic acid	172	C10H20O2	000334-48-5	62
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037442.D
Acq On : 09 Nov 2022 18:16
Operator : CG/JU
Sample : N5436-05 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
250FERRYBLVD-TFS-05

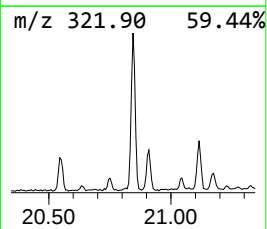
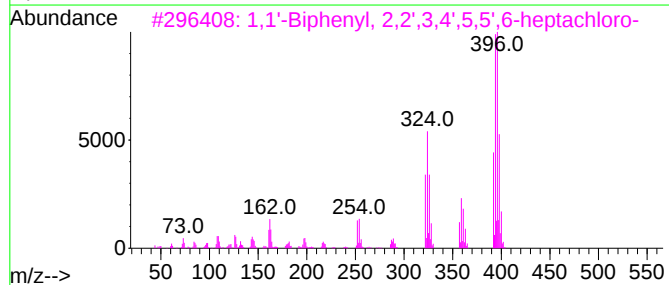
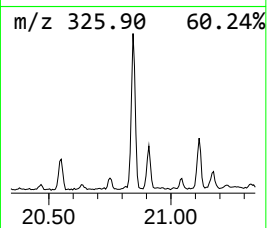
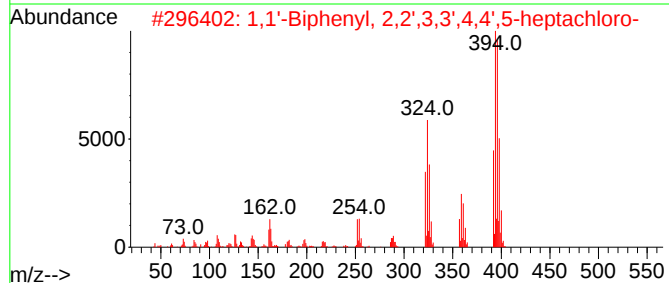
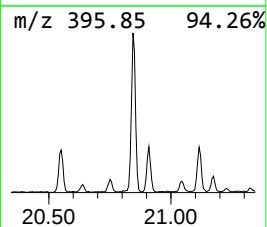
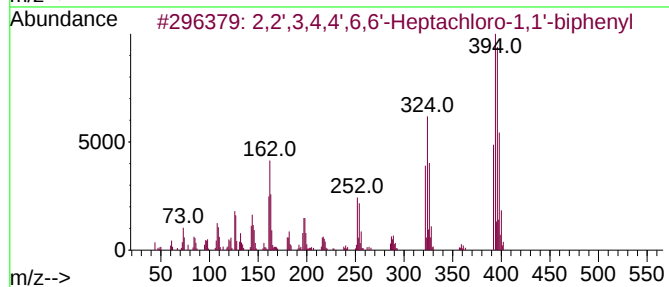
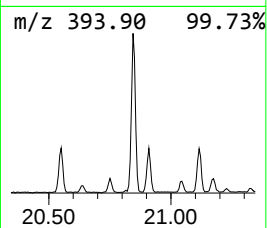
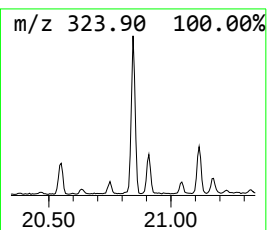
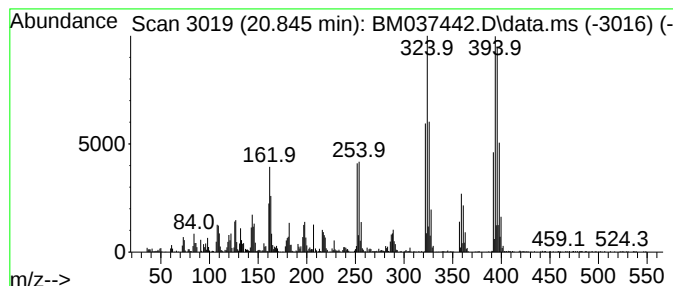
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2,2',3,4,4',6,6'-Heptachlor... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.845	3.03 ng	984559	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4,4',6,6'-Heptachloro-1,1...	392	C12H3Cl7	074472-48-3	99		
2	1,1'-Biphenyl, 2,2',3,3',4,4',5-...	392	C12H3Cl7	035065-30-6	99		
3	1,1'-Biphenyl, 2,2',3,4',5,5',6-...	392	C12H3Cl7	052663-68-0	99		
4	1,1'-Biphenyl, 2,2',3,3',4,4',6-...	392	C12H3Cl7	052663-71-5	99		
5	1,1'-Biphenyl, 2,2',3,3',5,6,6'-...	392	C12H3Cl7	052663-64-6	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037442.D
Acq On : 09 Nov 2022 18:16
Operator : CG/JU
Sample : N5436-05 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-05

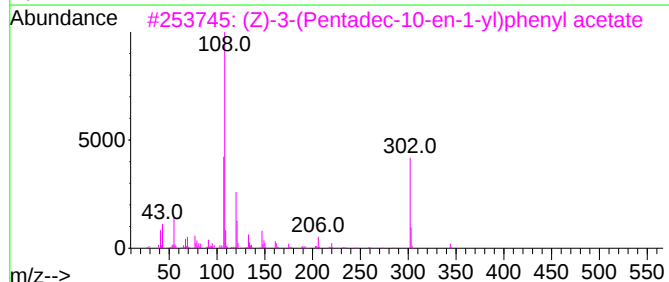
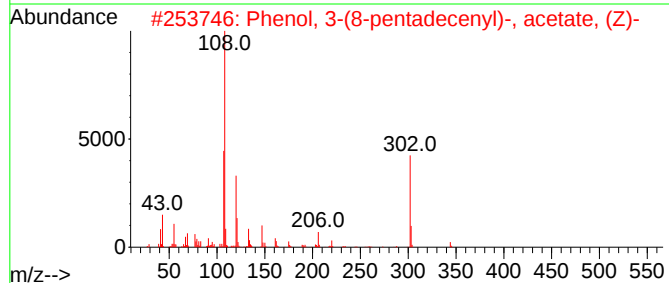
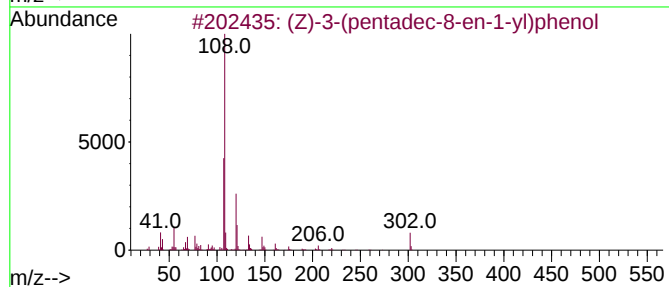
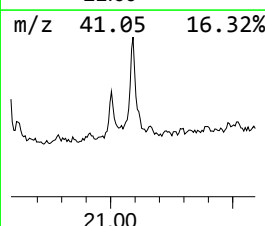
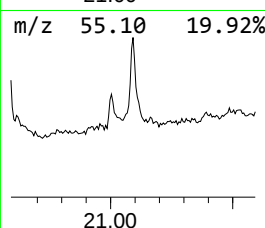
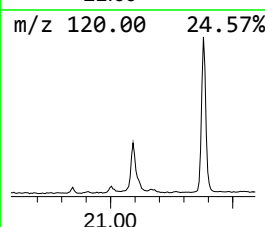
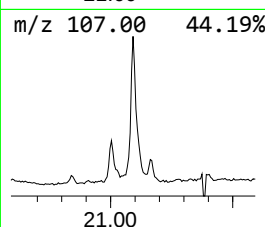
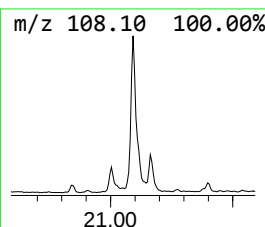
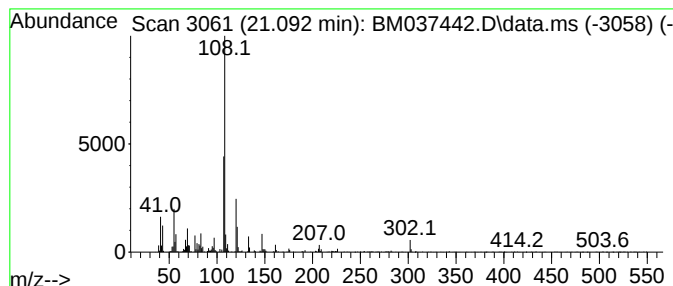
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 (Z)-3-(pentadec-8-en-1-yl)p... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.092	2.17 ng	704763	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(Z)-3-(pentadec-8-en-1-yl)phenol	302	C21H34O	000501-26-8	92
2			Phenol, 3-(8-pentadecenyl)-, ace...	344	C23H36O2	111047-34-8	83
3			(Z)-3-(Pentadec-10-en-1-yl)pheny...	344	C23H36O2	1000478-00-3	78
4			Acetic acid, 4-methylphenyl ester	150	C9H10O2	000140-39-6	50
5			Phenol, 3-pentadecyl-	304	C21H36O	000501-24-6	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037442.D
Acq On : 09 Nov 2022 18:16
Operator : CG/JU
Sample : N5436-05 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

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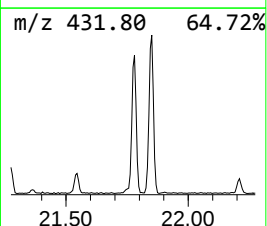
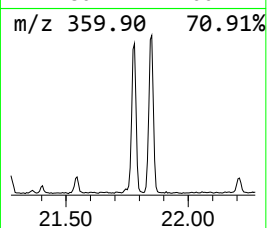
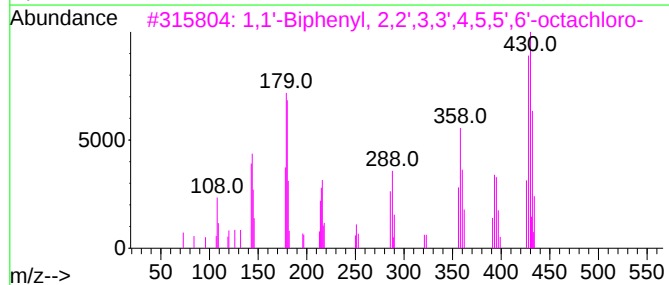
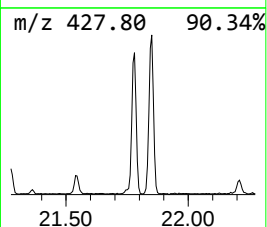
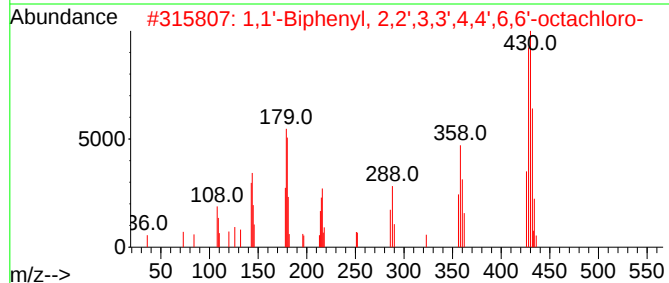
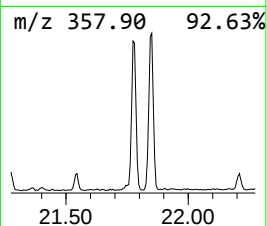
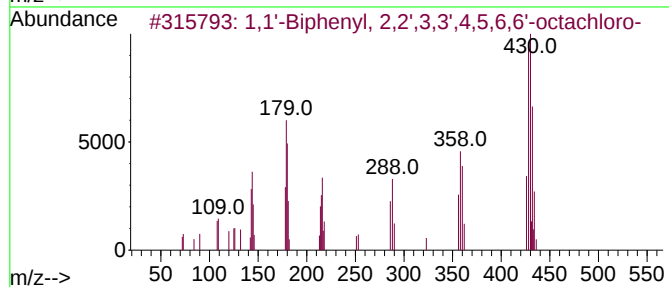
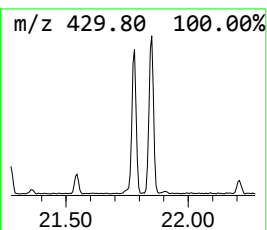
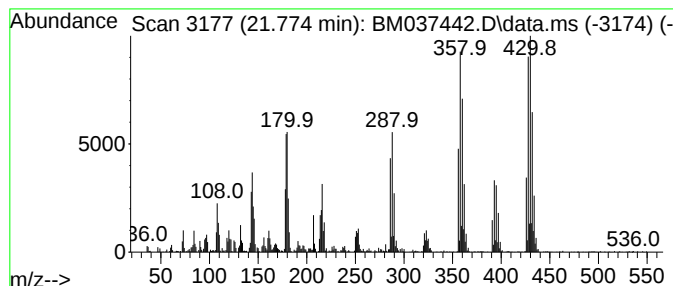
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.774	4.56 ng	1481870	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,3',4,5,6,6...	426	C12H2Cl8	052663-73-7	99		
2	1,1'-Biphenyl, 2,2',3,3',4,4',6...	426	C12H2Cl8	033091-17-7	99		
3	1,1'-Biphenyl, 2,2',3,3',4,5,5',...	426	C12H2Cl8	052663-75-9	99		
4	1,1'-Biphenyl, 2,2',3,3',4,4',5...	426	C12H2Cl8	042740-50-1	99		
5	1,1'-Biphenyl, 2,2',3,4,4',5,6,6...	426	C12H2Cl8	074472-52-9	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037442.D
Acq On : 09 Nov 2022 18:16
Operator : CG/JU
Sample : N5436-05 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

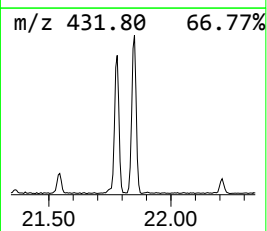
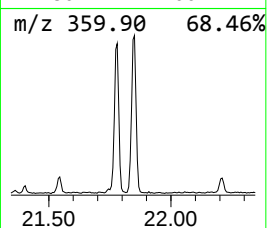
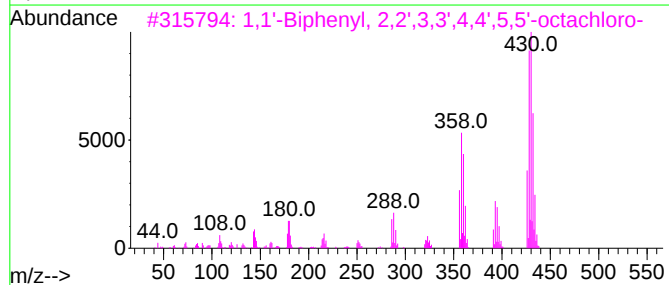
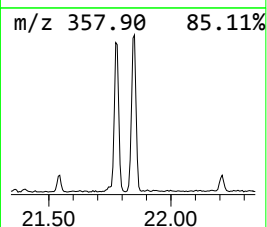
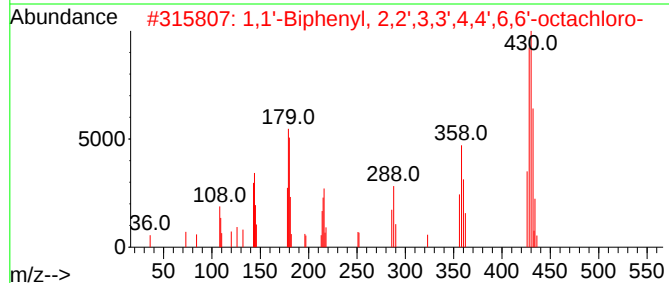
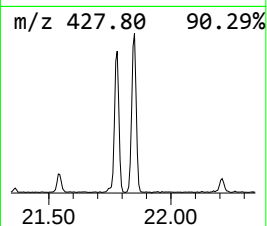
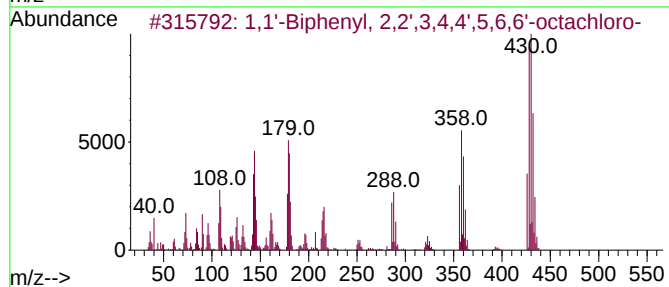
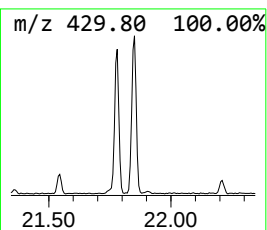
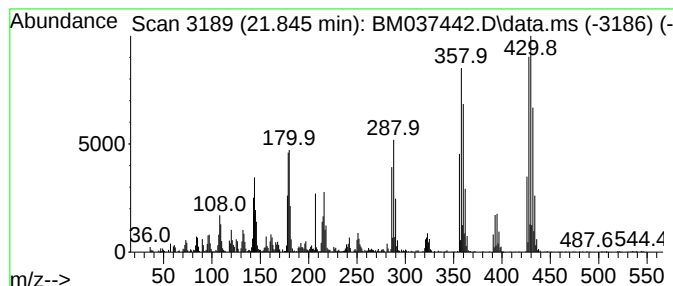
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1,1'-Biphenyl, 2,2',3,4,4',... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.845	4.32 ng	1402570	Chrysene-d12	21.380	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,4,4',5,6,6...	426	C12H2Cl8	074472-52-9	99
2	1,1'-Biphenyl, 2,2',3,3',4,4',6...	426	C12H2Cl8	033091-17-7	99
3	1,1'-Biphenyl, 2,2',3,3',4,4',5...	426	C12H2Cl8	035694-08-7	99
4	1,1'-Biphenyl, 2,2',3,3',4,4',5...	426	C12H2Cl8	042740-50-1	99
5	1,1'-Biphenyl, 2,2',3,3',4,4',5...	426	C12H2Cl8	052663-78-2	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037442.D
Acq On : 09 Nov 2022 18:16
Operator : CG/JU
Sample : N5436-05 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

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ClientSampleId :
250FERRYBLVD-TFS-05

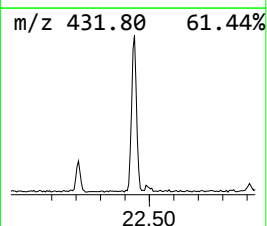
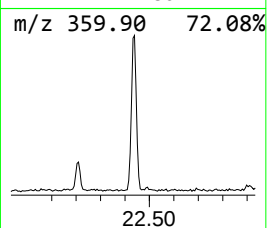
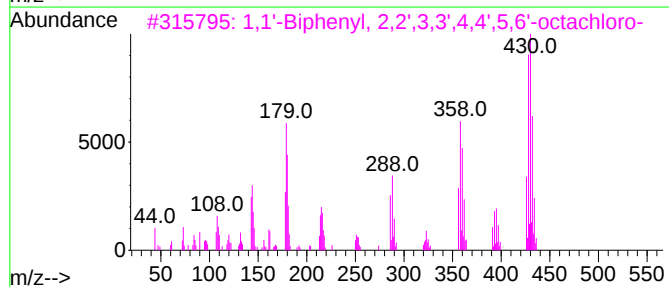
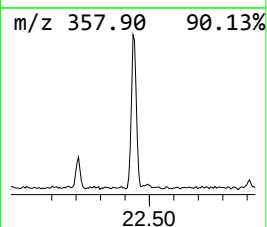
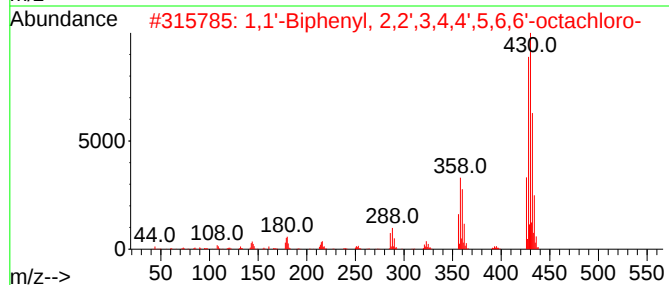
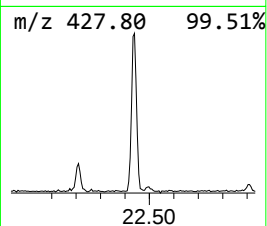
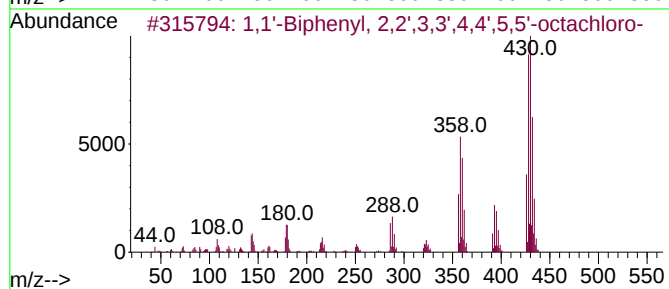
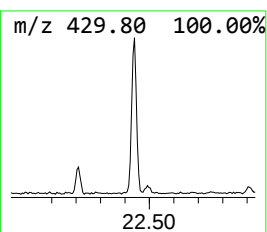
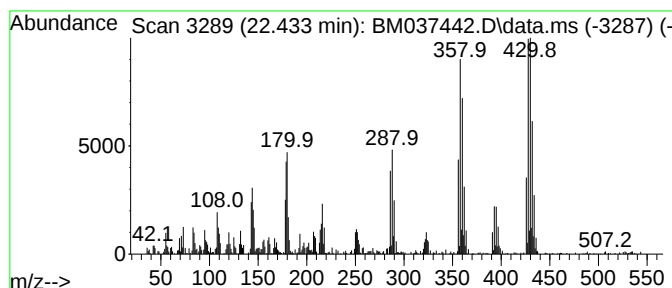
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.433	2.40 ng	777785	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,3',4,4',5,...	426	C12H2Cl8	035694-08-7	99		
2	1,1'-Biphenyl, 2,2',3,4,4',5,6,6...	426	C12H2Cl8	074472-52-9	99		
3	1,1'-Biphenyl, 2,2',3,3',4,4',5,...	426	C12H2Cl8	042740-50-1	99		
4	1,1'-Biphenyl, 2,2',3,3',4,5,5',...	426	C12H2Cl8	052663-75-9	99		
5	1,1'-Biphenyl, 2,2',3,3',4,4',5,...	426	C12H2Cl8	052663-78-2	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037442.D
Acq On : 09 Nov 2022 18:16
Operator : CG/JU
Sample : N5436-05 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

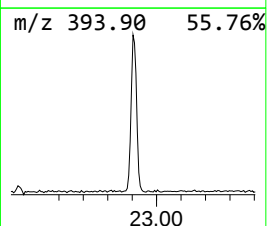
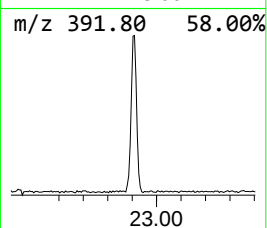
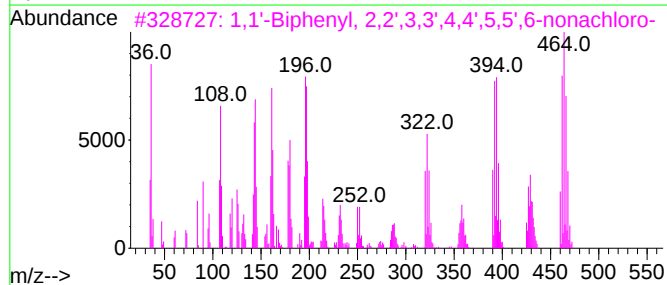
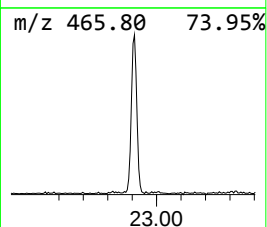
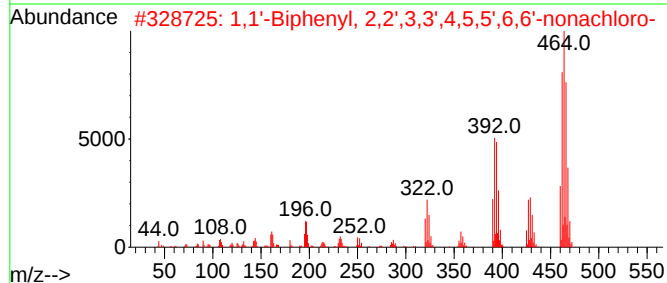
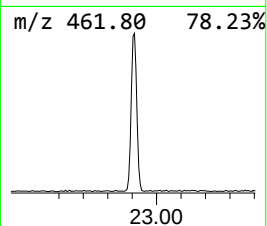
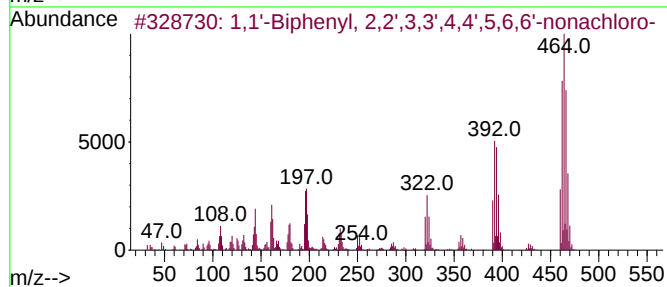
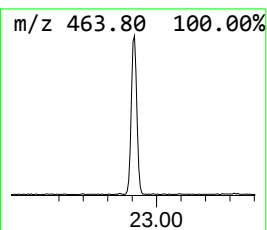
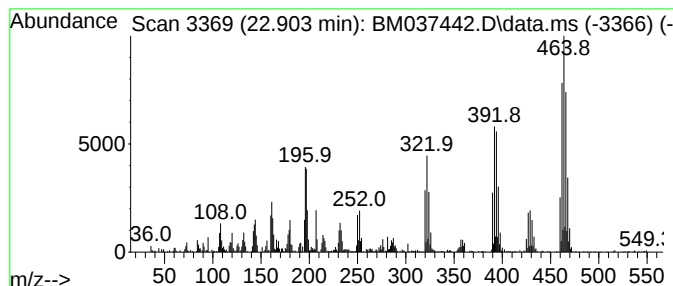
Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-05

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.904	5.99 ng	1188890	Perylene-d12	23.721
Hit#	of 5	Tentative ID	MW MolForm	CAS# Qual
1	1,1'-Biphenyl, 2,2',3,3',4,4',5,...	460 C12HC19	052663-79-3	99
2	1,1'-Biphenyl, 2,2',3,3',4,5,5',...	460 C12HC19	052663-77-1	99
3	1,1'-Biphenyl, 2,2',3,3',4,4',5,...	460 C12HC19	040186-72-9	83
4	Spiro[benzofuran-2(3H),1'-[2,5]c...	462 C14H9Br3O3	052498-93-8	35
5	13,27-Cyclours-11-en-3-ol, acetate	466 C32H50O2	1000189-80-7	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037442.D
Acq On : 09 Nov 2022 18:16
Operator : CG/JU
Sample : N5436-05 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-05

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.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
2-Pentanone, 4-...	4.934	2.3	ng	187134	1	7.822	1653870	20.0
o-Terphenyl	17.851	12.9	ng	2404200	4	17.204	3722540	20.0
n-Hexadecanoic ...	18.098	8.8	ng	1637670	4	17.204	3722540	20.0
Phenol, 2-[(4-h...	18.604	4.7	ng	880391	4	17.204	3722540	20.0
Phenol, 4,4'-me...	19.015	5.6	ng	1046580	4	17.204	3722540	20.0
Octadecanoic acid	19.374	2.8	ng	893902	5	21.380	6493290	20.0
2,2',3,4,4',6,6...	20.845	3.0	ng	984559	5	21.380	6493290	20.0
(Z)-3-(pentadec...	21.092	2.2	ng	704763	5	21.380	6493290	20.0
1,1'-Biphenyl, ...	21.774	4.6	ng	1481870	5	21.380	6493290	20.0
1,1'-Biphenyl, ...	21.845	4.3	ng	1402570	5	21.380	6493290	20.0
1,1'-Biphenyl, ...	22.433	2.4	ng	777785	5	21.380	6493290	20.0
1,1'-Biphenyl, ...	22.904	6.0	ng	1188890	6	23.721	3970400	20.0