

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
 Data File : BM037441.D
 Acq On : 09 Nov 2022 17:40
 Operator : CG/JU
 Sample : N5436-10 10X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 250FERRYBLVD-TFS-10

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: BM037441.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	309	314	322	rVB	152746	205876	1.63%	0.317%
2	5.398	387	393	398	rBV	486284	708233	5.61%	1.091%
3	7.004	661	666	679	rVB	427352	710303	5.62%	1.095%
4	7.822	799	805	818	rVB	1113578	1709486	13.54%	2.635%
5	8.610	934	939	950	rBV2	58989	132456	1.05%	0.204%
6	8.992	997	1004	1009	rBV	270273	464191	3.68%	0.715%
7	9.810	1135	1143	1154	rBV	126132	273777	2.17%	0.422%
8	10.045	1176	1183	1192	rBV2	46368	134244	1.06%	0.207%
9	10.622	1269	1281	1286	rBV	1403138	2308951	18.28%	3.558%
10	10.669	1286	1289	1295	rVB	168283	271697	2.15%	0.419%
11	13.086	1692	1700	1708	rBV	722419	1051161	8.32%	1.620%
12	14.186	1882	1887	1897	rVB	90024	146141	1.16%	0.225%
13	14.463	1925	1934	1940	rBV	2108130	3075297	24.35%	4.740%
14	14.521	1940	1944	1952	rVB	178820	282739	2.24%	0.436%
15	14.857	1996	2001	2010	rBV	98080	179129	1.42%	0.276%
16	15.504	2106	2111	2122	rBV	155536	249745	1.98%	0.385%
17	15.710	2142	2146	2151	rBV4	90320	134747	1.07%	0.208%
18	15.951	2182	2187	2193	rBV	618054	856828	6.79%	1.321%
19	16.180	2219	2226	2232	rBV6	75527	196473	1.56%	0.303%
20	17.204	2394	2400	2404	rBV	2599418	3603799	28.54%	5.554%
21	17.245	2404	2407	2415	rVV	673508	947843	7.51%	1.461%
22	17.339	2419	2423	2427	rVV	166141	227945	1.81%	0.351%
23	17.851	2505	2510	2517	rVB	1846631	2305058	18.25%	3.552%
24	18.098	2546	2552	2563	rBV2	732794	1486682	11.77%	2.291%
25	18.274	2577	2582	2586	rBV2	238905	389759	3.09%	0.601%
26	18.392	2597	2602	2606	rBV2	201060	327006	2.59%	0.504%
27	18.603	2631	2638	2651	rVB3	641958	1252108	9.92%	1.930%
28	18.980	2699	2702	2704	rBV	223021	287297	2.28%	0.443%
29	19.015	2704	2708	2718	rVB	715339	1126671	8.92%	1.736%
30	19.262	2745	2750	2755	rBV	1550399	2200561	17.43%	3.391%
31	19.309	2755	2758	2763	rVB2	342689	452148	3.58%	0.697%
32	19.374	2766	2769	2773	rBV	498058	697492	5.52%	1.075%
33	19.621	2803	2811	2820	rBV	9728002	12627977	100.00%	19.462%
34	19.827	2842	2846	2848	rBV	1196381	1375470	10.89%	2.120%
35	19.862	2848	2852	2858	rVB	5503607	6189369	49.01%	9.539%
36	20.268	2917	2921	2926	rBV2	363994	532462	4.22%	0.821%

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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

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37	20.550	2966	2969	2972	rBV3	249887	316049	2.50%	0.487%
38	20.586	2972	2975	2979	rVV	295939	318967	2.53%	0.492%
39	20.844	3016	3019	3025	rBV2	753437	843517	6.68%	1.300%
40	21.003	3043	3046	3049	rBV	482103	611105	4.84%	0.942%
41	21.092	3058	3061	3064	rBV	597286	708134	5.61%	1.091%
42	21.380	3105	3110	3122	rVB2	3441628	6729306	53.29%	10.371%
43	21.774	3174	3177	3184	rVB	921232	1177523	9.32%	1.815%
44	21.844	3186	3189	3192	rBV	907553	1165385	9.23%	1.796%
45	23.721	3503	3508	3515	rVB2	2106598	3895322	30.85%	6.003%

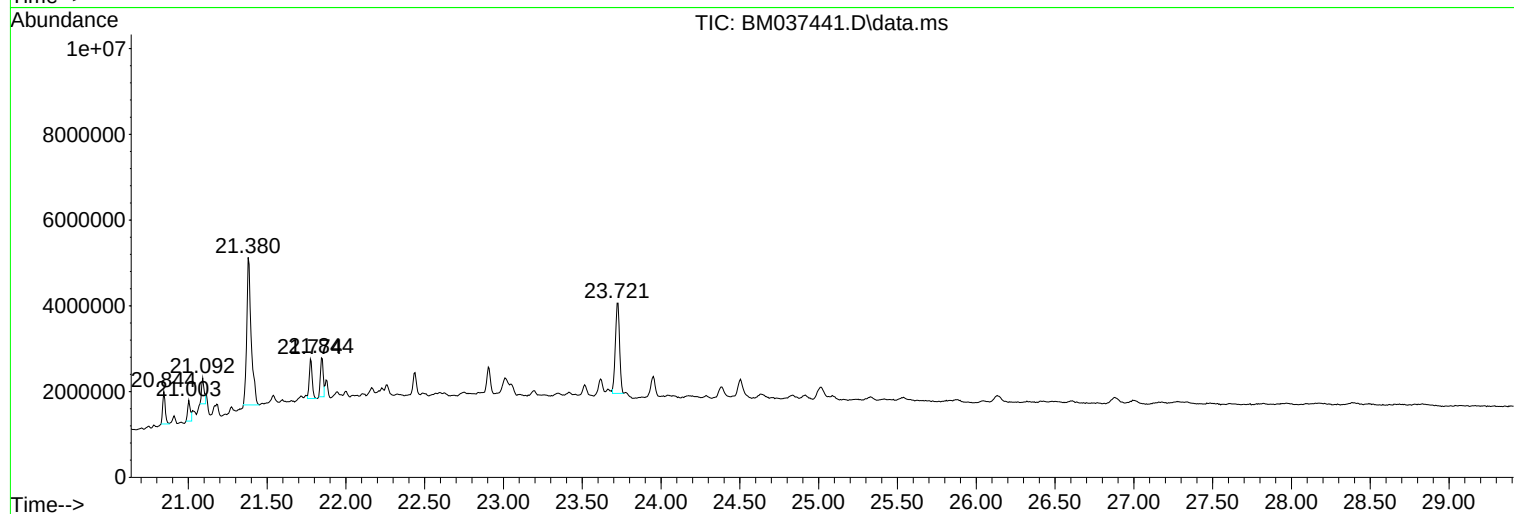
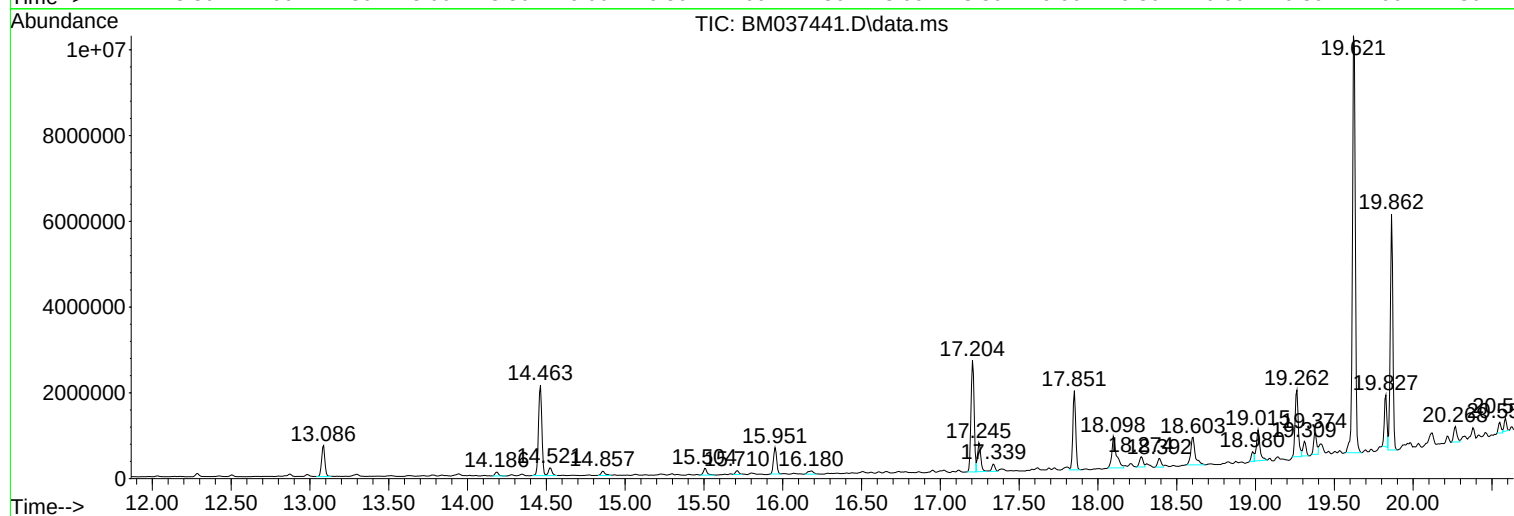
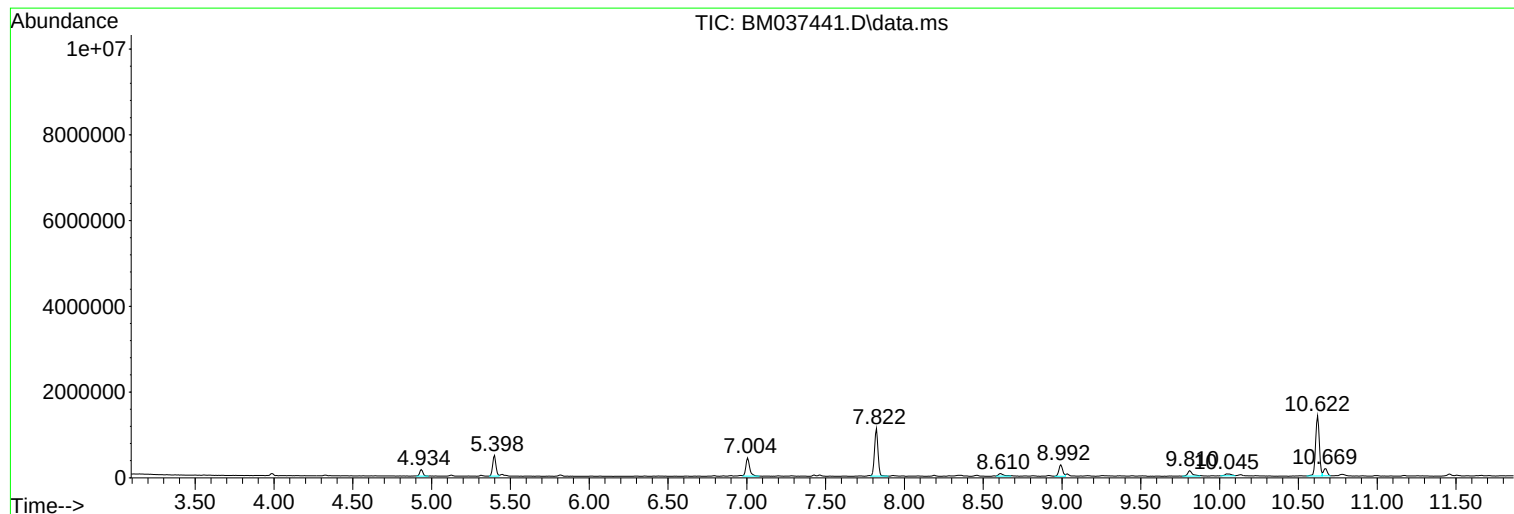
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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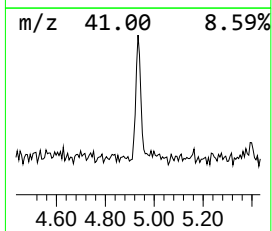
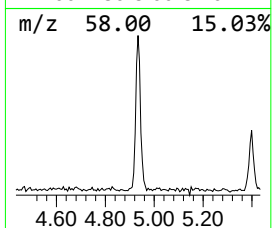
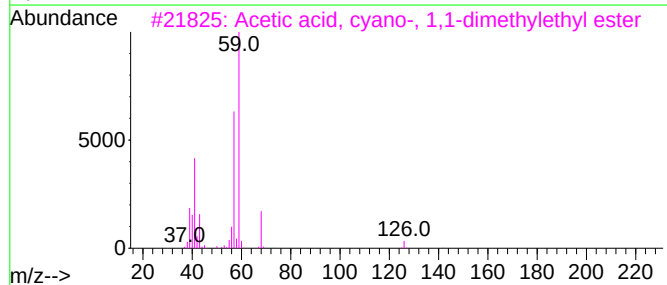
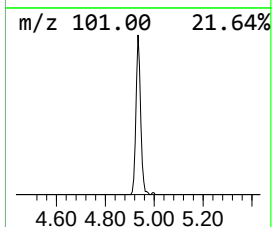
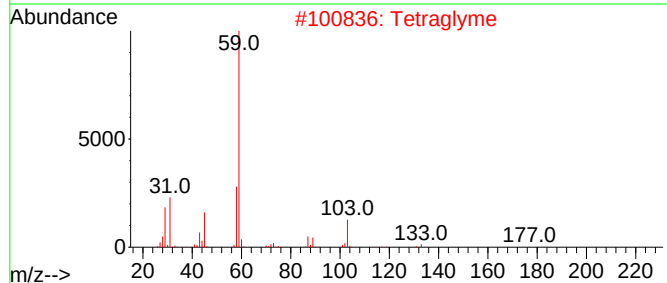
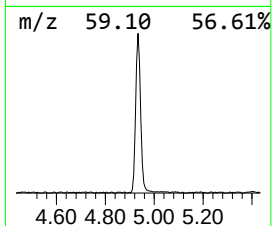
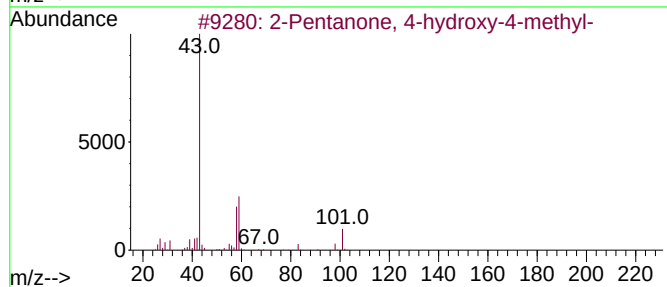
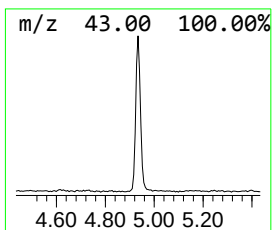
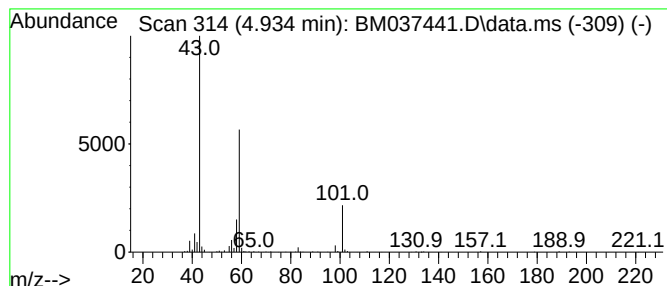
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.934	2.41 ng	205876	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	64		
2	Tetraglyme	222	C10H22O5	000143-24-8	25		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23		
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		



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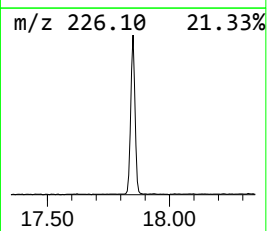
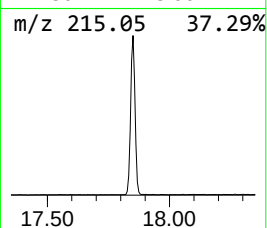
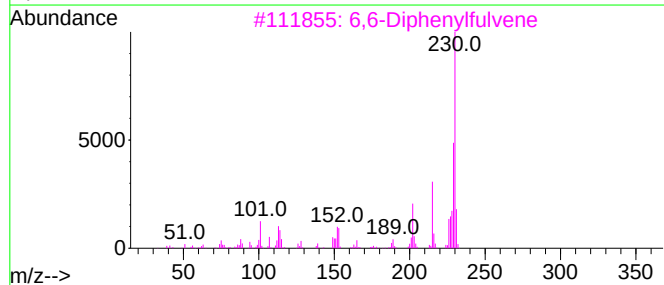
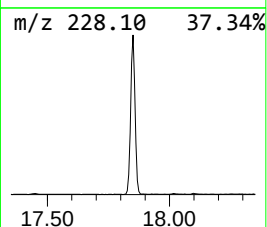
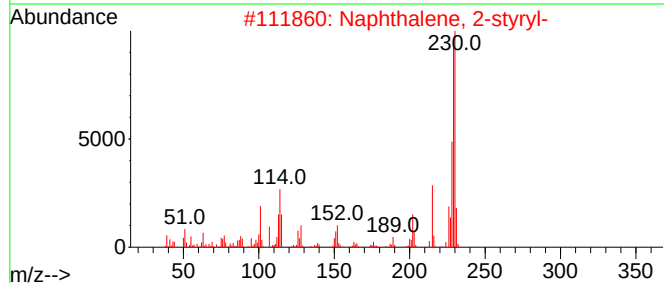
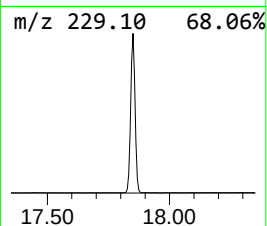
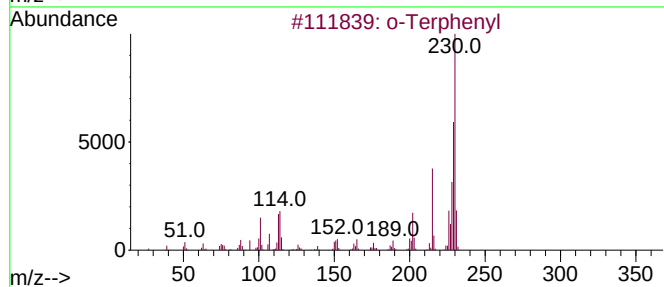
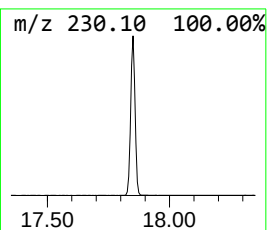
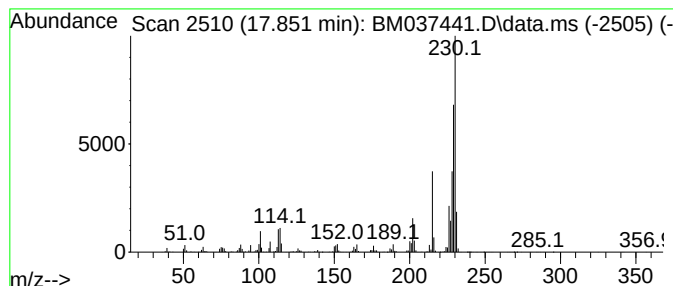
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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	12.79 ng	2305060	Phenanthrene-d10	17.203

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Terphenyl	230	C18H14	000084-15-1	99
2			Naphthalene, 2-styryl-	230	C18H14	002039-70-5	93
3			6,6-Diphenylfulvene	230	C18H14	002175-90-8	93
4			Pyrene, 1,3-dimethyl-	230	C18H14	064401-21-4	86
5			Naphthacene, 5,12-dihydro-	230	C18H14	000959-02-4	62



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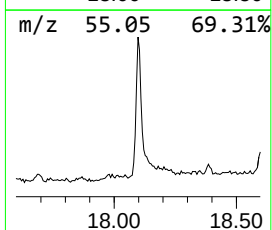
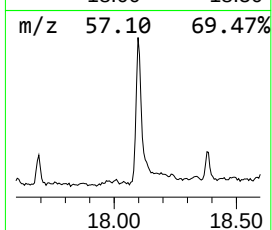
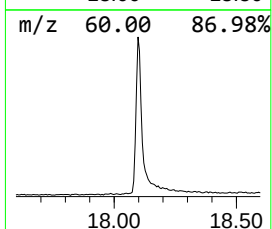
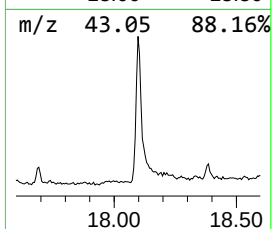
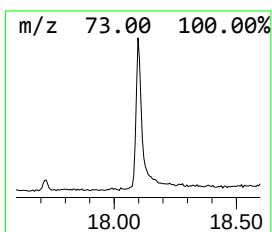
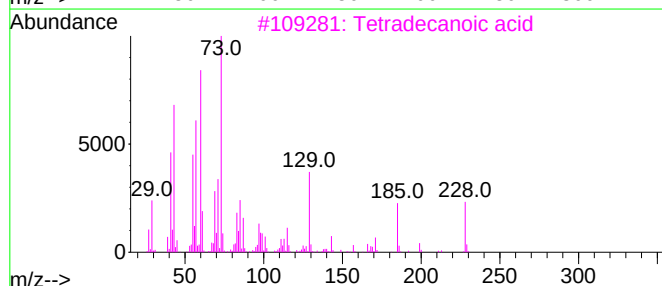
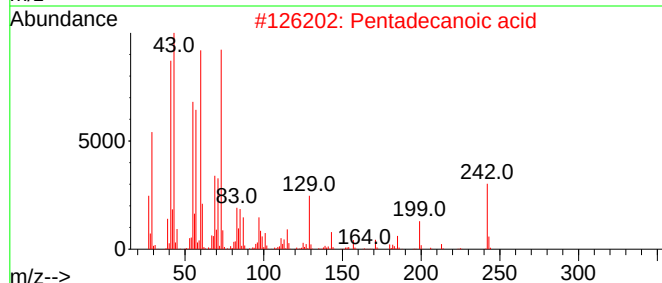
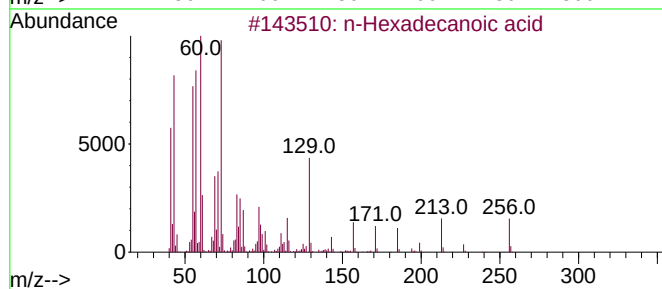
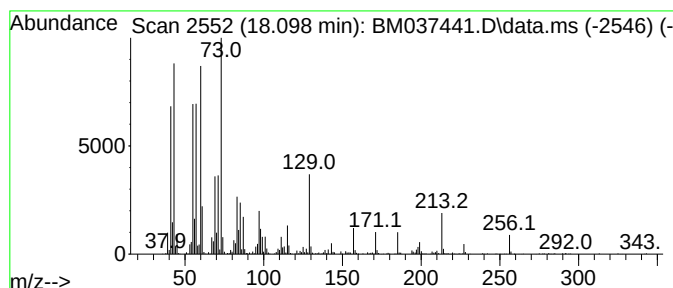
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.098	8.25 ng	1486680	Phenanthrene-d10	17.203

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	91
4			Octadecanoic acid	284	C18H36O2	000057-11-4	87
5			Tridecanoic acid	214	C13H26O2	000638-53-9	81



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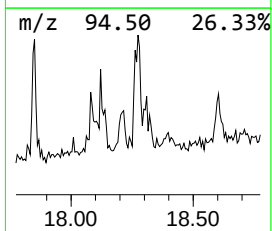
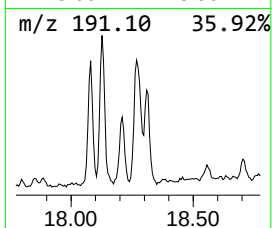
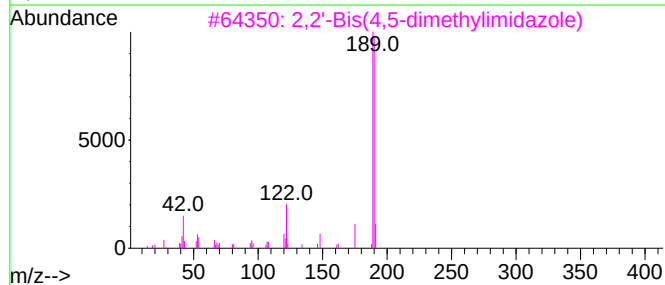
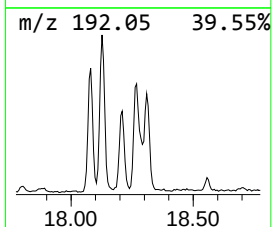
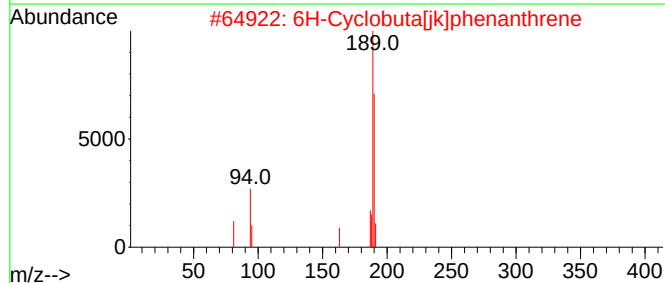
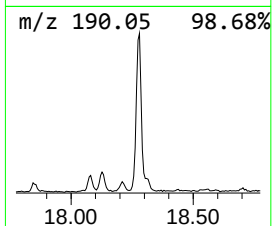
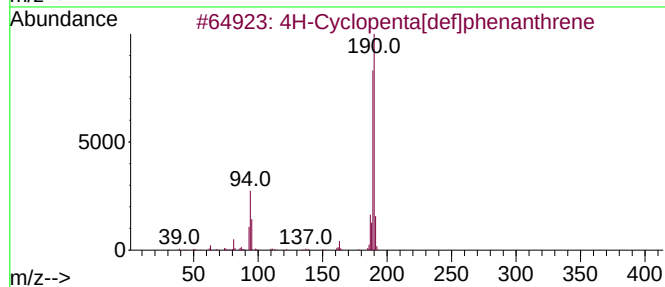
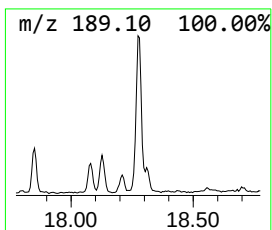
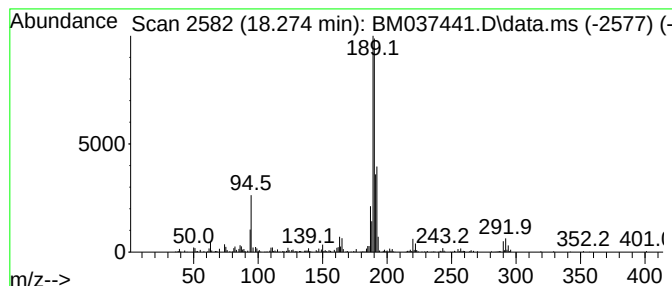
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.274	2.16 ng	389759	Phenanthrene-d10	17.203

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	59
3			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	47
4			Methyl diselenide	190	C2H6Se2	007101-31-7	43
5			Pyrimidine, 6-oxo-4-(3-fluorophe...	190	C10H7FN2O	085979-55-1	27



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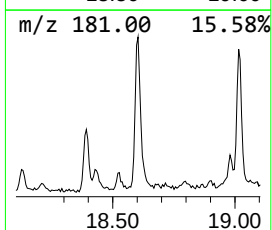
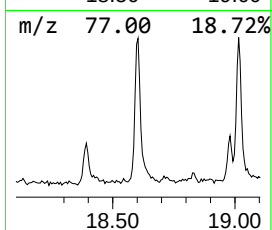
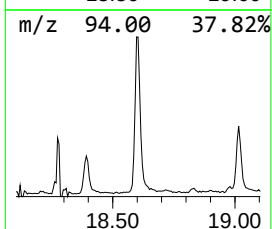
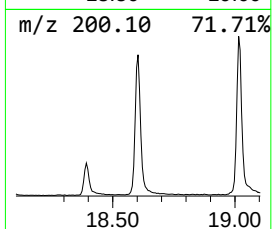
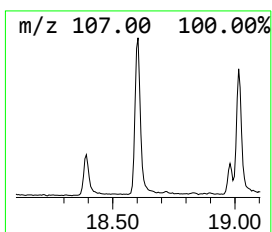
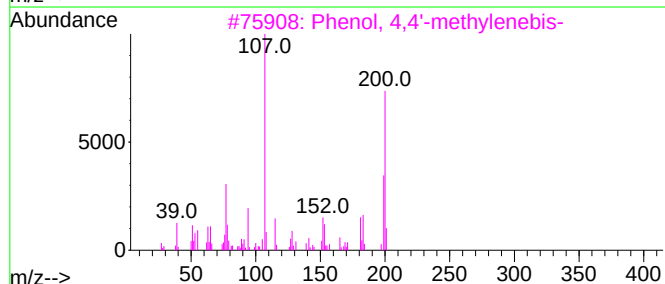
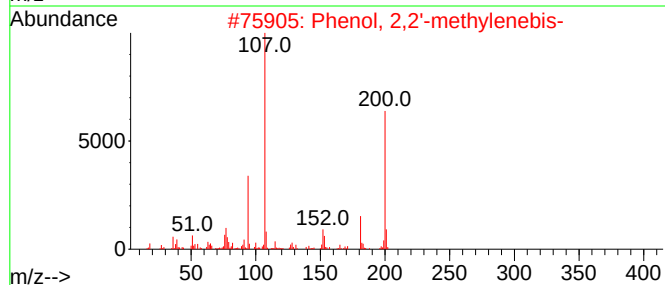
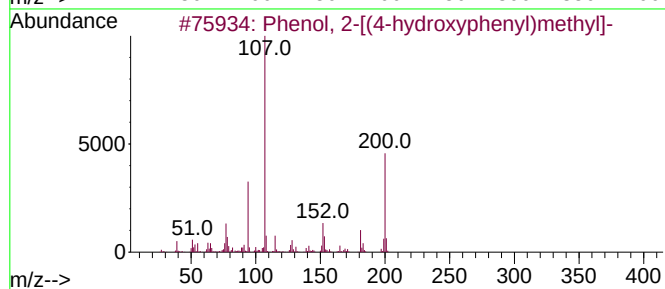
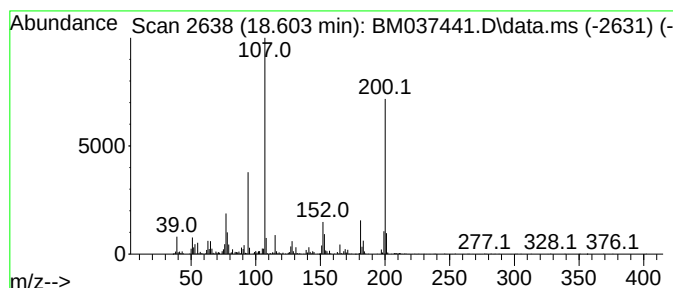
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Phenol, 2-[(4-hydroxyphenyl)... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.603	6.95 ng	1252110	Phenanthrene-d10	17.203

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	97
2			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	94
3			Phenol, 4,4'-methylenebis-	200	C13H12O2	000620-92-8	72
4			1,3-Benzenediol, 4-(phenylmethyl)-	200	C13H12O2	002284-30-2	42
5			4-(2-Methoxyethyl)phenol	152	C9H12O2	056718-71-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037441.D
Acq On : 09 Nov 2022 17:40
Operator : CG/JU
Sample : N5436-10 10X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-10

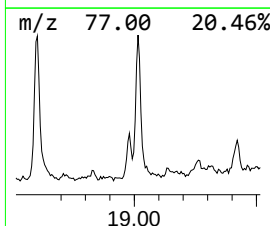
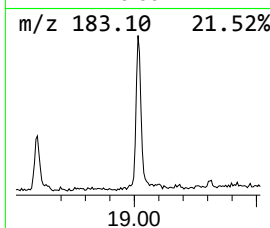
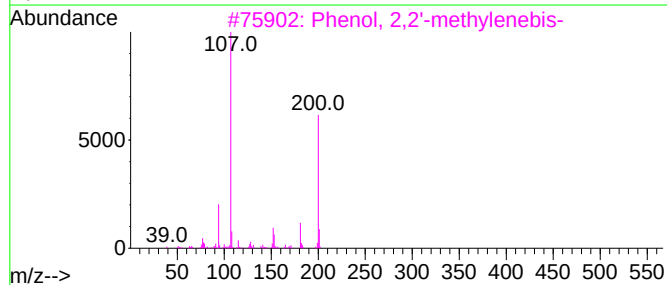
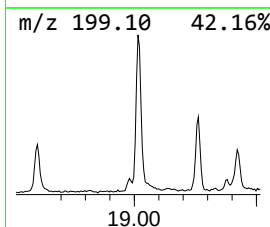
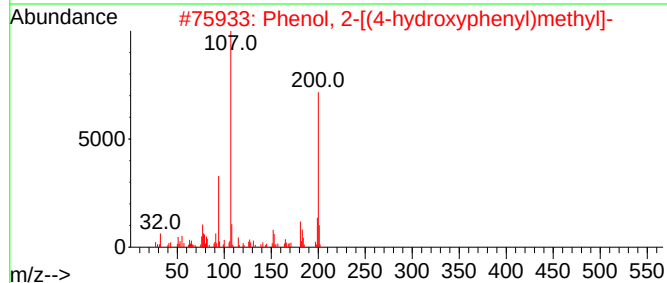
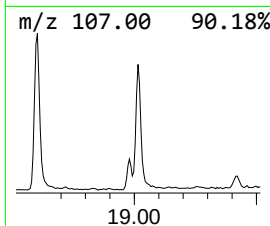
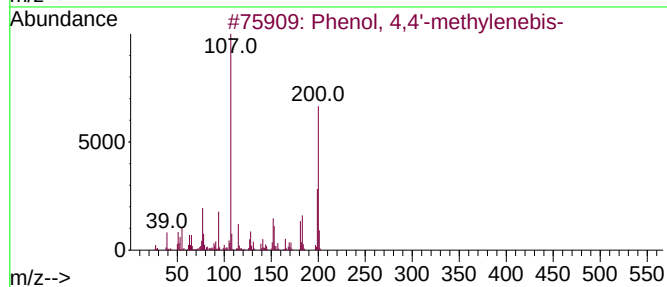
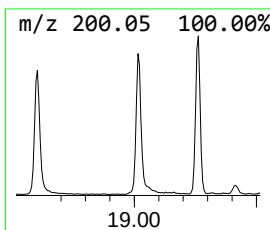
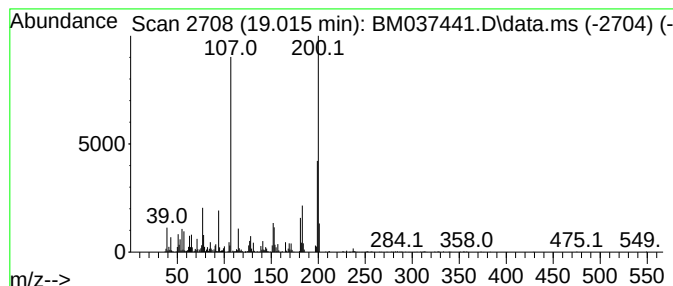
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phenol, 4,4'-methylenebis- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.015	6.25 ng	1126670	Phenanthrene-d10	17.203

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4,4'-methylenebis-	200	C13H12O2	000620-92-8	97
2			Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	90
3			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	81
4			2,4'-Dihydroxydiphenylmethane, d...	284	C17H16O4	959033-36-4	64
5			1,3-Benzenediol, 4-(phenylmethyl)-	200	C13H12O2	002284-30-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037441.D
Acq On : 09 Nov 2022 17:40
Operator : CG/JU
Sample : N5436-10 10X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-10

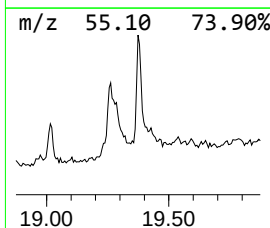
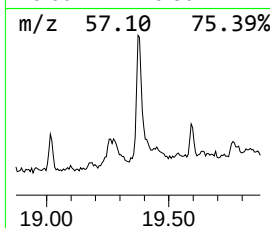
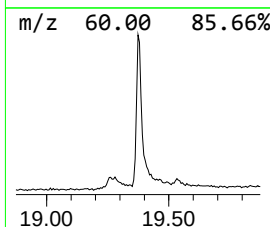
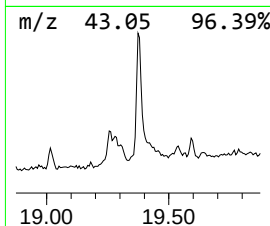
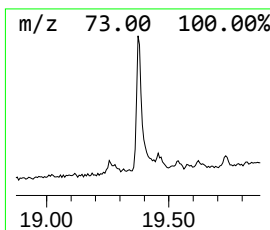
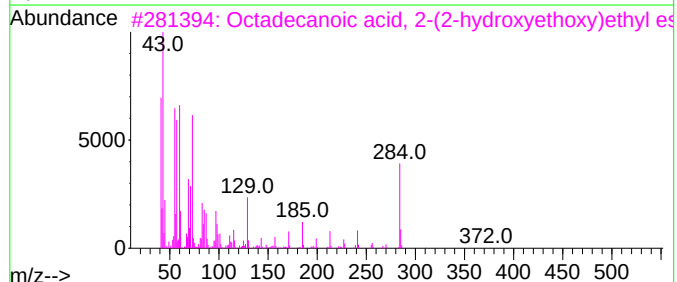
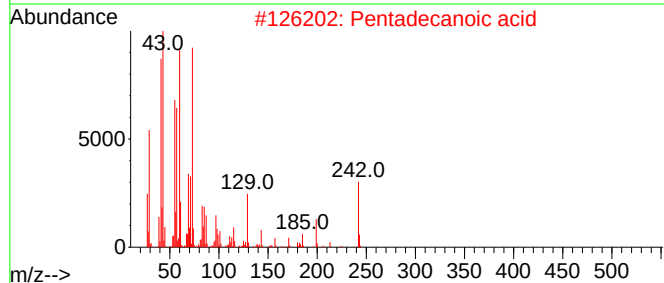
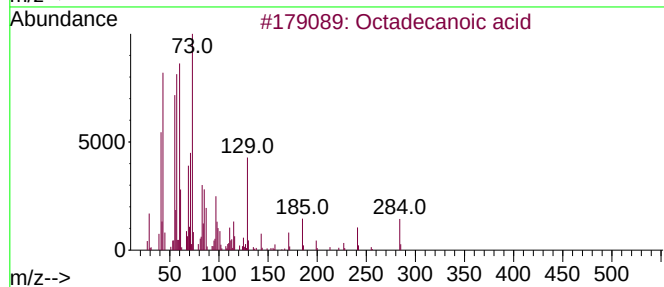
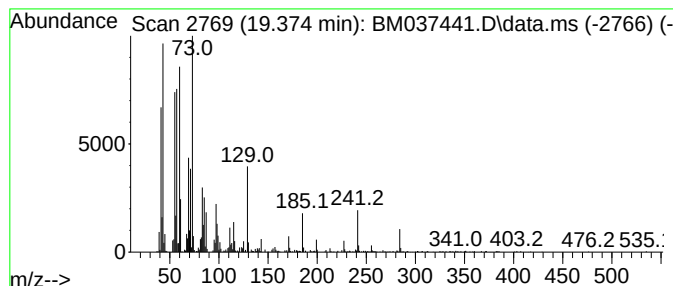
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.374	2.07 ng	697492	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	91
3			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	83
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
5			n-Decanoic acid	172	C10H20O2	000334-48-5	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037441.D
Acq On : 09 Nov 2022 17:40
Operator : CG/JU
Sample : N5436-10 10X
Misc :
ALS Vial : 10 Sample Multiplier: 1

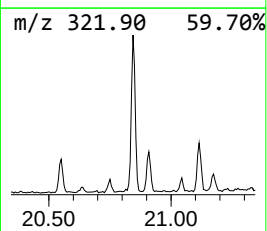
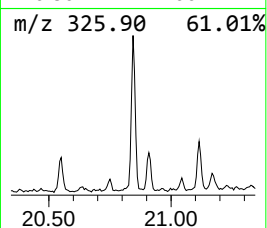
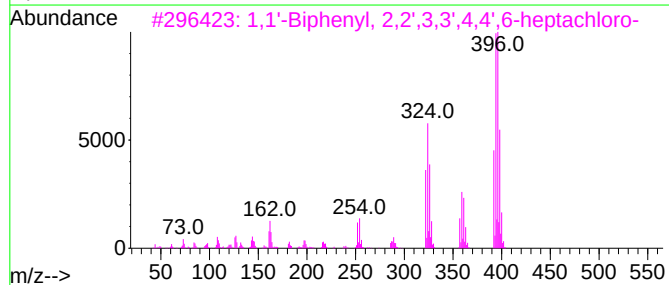
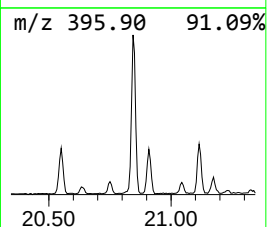
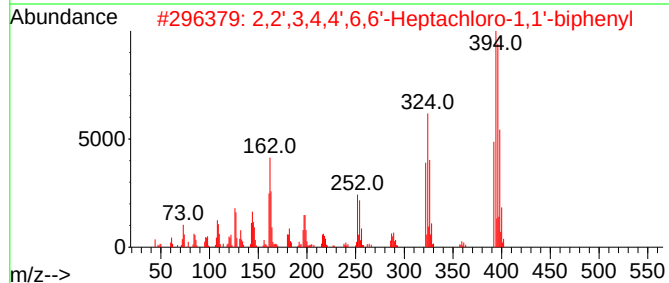
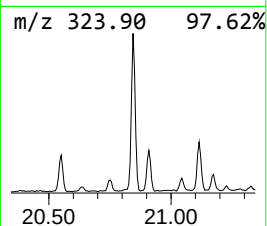
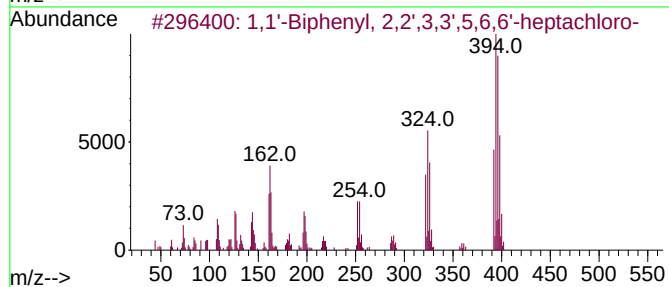
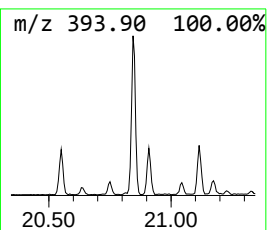
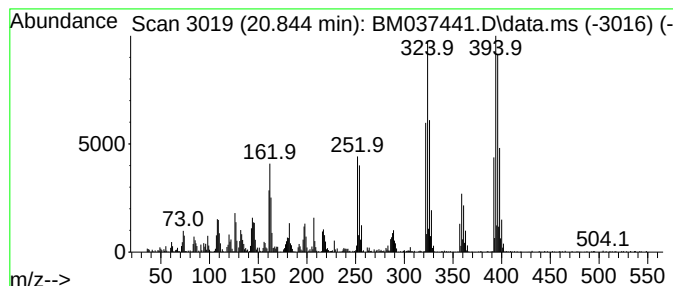
Instrument :
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ClientSampleId :
250FERRYBLVD-TFS-10

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.845	2.51 ng	843517	Chrysene-d12		21.380	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,3',5,6,6'-...		392	C12H3Cl7	052663-64-6	99
2	2,2',3,4,4',6,6'-Heptachloro-1,1...		392	C12H3Cl7	074472-48-3	99
3	1,1'-Biphenyl, 2,2',3,3',4,4',6-...		392	C12H3Cl7	052663-71-5	99
4	1,1'-Biphenyl, 2,2',3,3',4,4',5-...		392	C12H3Cl7	035065-30-6	99
5	1,1'-Biphenyl, 2,2',3,4',5,5',6-...		392	C12H3Cl7	052663-68-0	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037441.D
Acq On : 09 Nov 2022 17:40
Operator : CG/JU
Sample : N5436-10 10X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
250FERRYBLVD-TFS-10

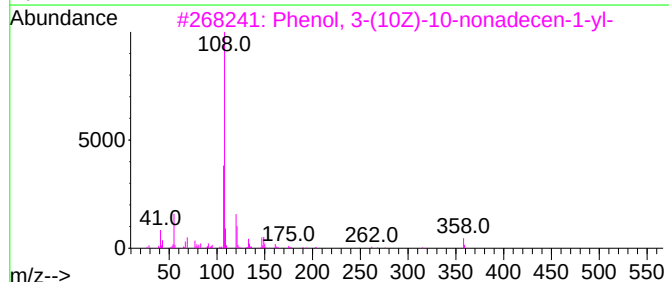
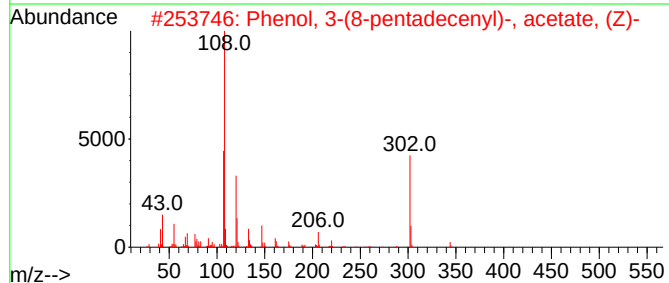
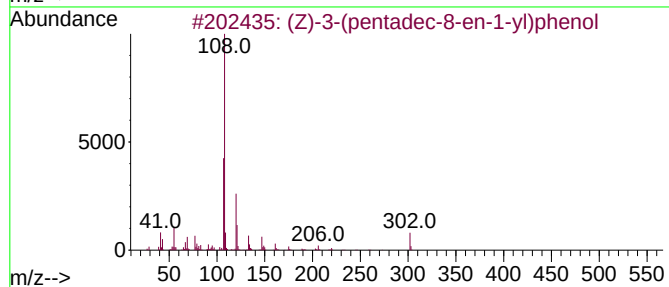
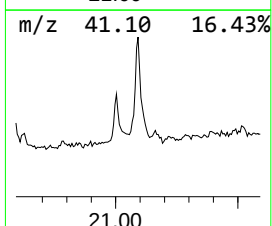
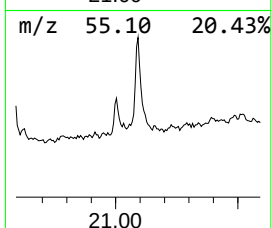
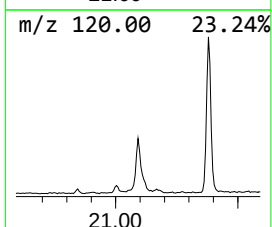
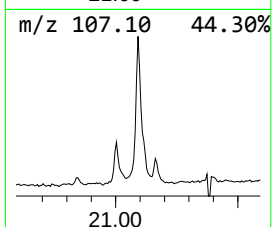
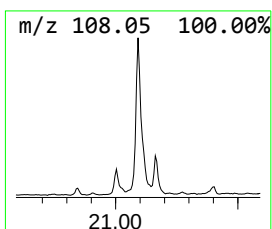
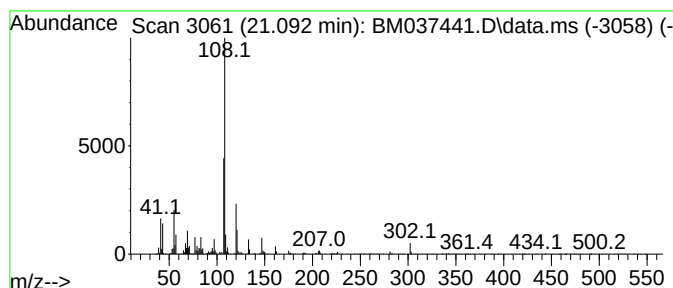
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 (Z)-3-(pentadec-8-en-1-yl)p... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.092	2.10 ng	708134	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	96
2		Phenol, 3-(8-pentadecenyl)-, ace...	344	C23H36O2	111047-34-8	83
3		Phenol, 3-(10Z)-10-nonadecen-1-yl-	358	C25H42O	936109-24-9	80
4		(Z)-3-(Pentadec-10-en-1-yl)pheny...	344	C23H36O2	1000478-00-3	78
5		Acetic acid, 4-methylphenyl ester	150	C9H10O2	000140-39-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037441.D
Acq On : 09 Nov 2022 17:40
Operator : CG/JU
Sample : N5436-10 10X
Misc :
ALS Vial : 10 Sample Multiplier: 1

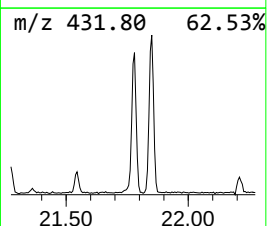
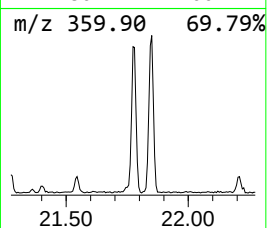
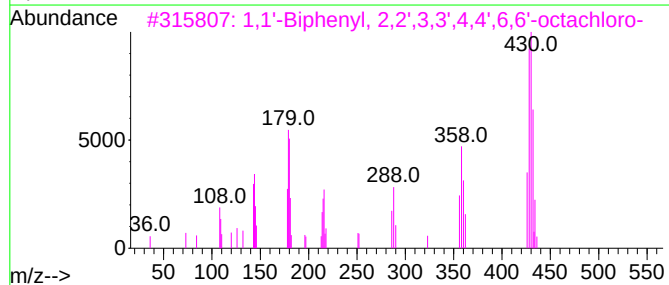
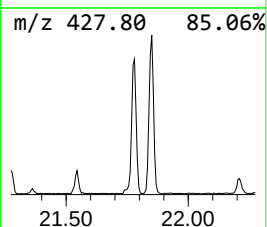
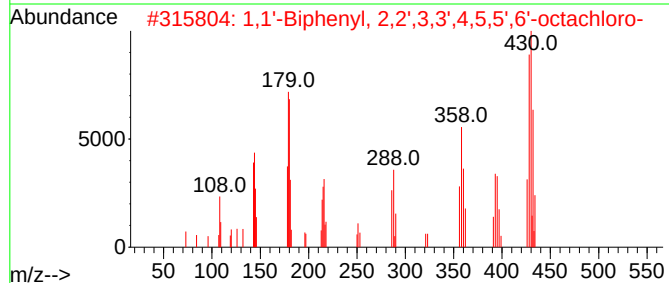
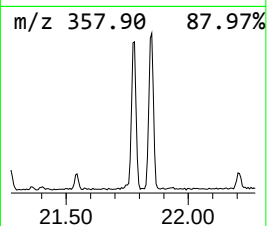
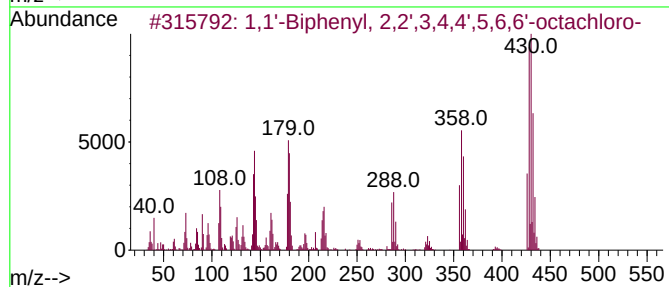
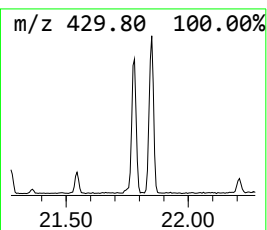
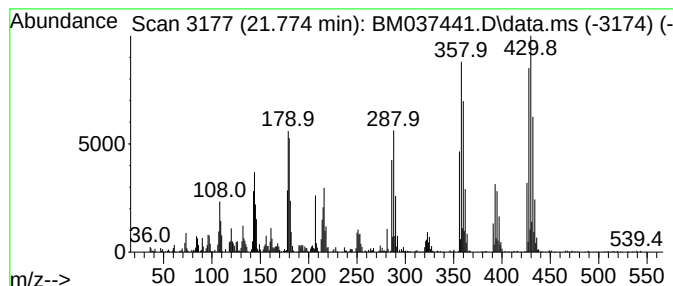
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1,1'-Biphenyl, 2,2',3,4,4',... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.774	3.50 ng	1177520	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',4,5,5',...	426	C12H2Cl8	052663-75-9	99
3	1,1'-Biphenyl, 2,2',3,3',4,4',6,...	426	C12H2Cl8	033091-17-7	99
4	1,1'-Biphenyl, 2,2',3,3',4,5,5',...	426	C12H2Cl8	068194-17-2	99
5	1,1'-Biphenyl, 2,2',3,3',4,5,6,6...	426	C12H2Cl8	052663-73-7	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
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Acq On : 09 Nov 2022 17:40
Operator : CG/JU
Sample : N5436-10 10X
Misc :
ALS Vial : 10 Sample Multiplier: 1

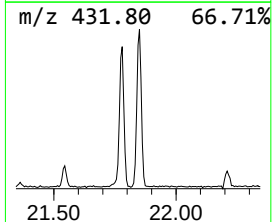
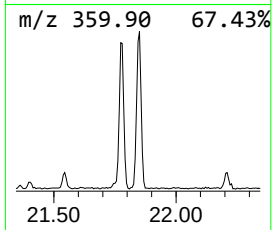
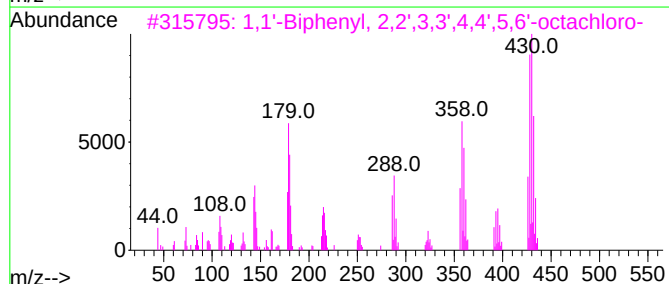
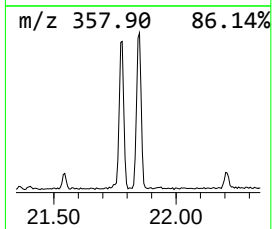
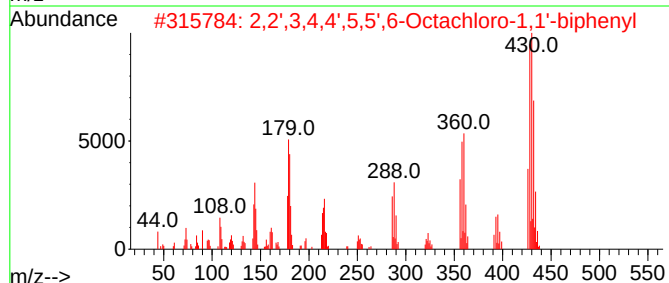
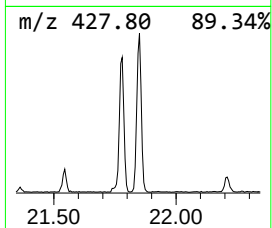
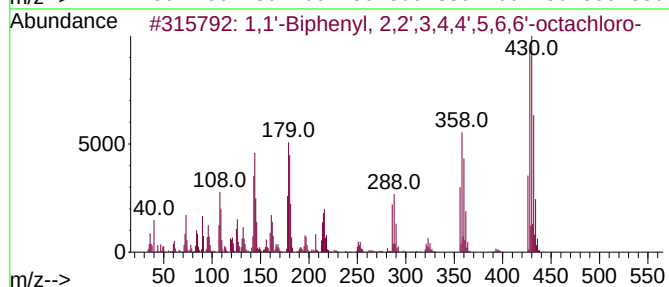
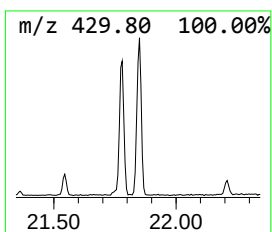
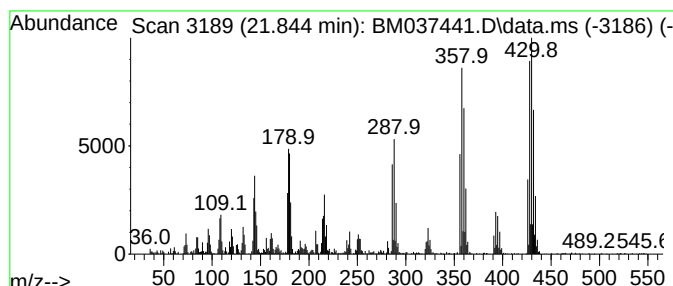
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.844	3.46 ng	1165390	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	2,2',3,4,4',5,5',6-Octachloro-1,...		426	C12H2Cl8	052663-76-0	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',4,4',5,...		426	C12H2Cl8	035694-08-7	99
5	1,1'-Biphenyl, 2,2',3,3',4,4',6,...		426	C12H2Cl8	033091-17-7	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037441.D
Acq On : 09 Nov 2022 17:40
Operator : CG/JU
Sample : N5436-10 10X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-10

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	2.4	ng	205876	1	7.822	1709490	20.0
o-Terphenyl	17.851	12.8	ng	2305060	4	17.203	3603800	20.0
n-Hexadecanoic ...	18.098	8.3	ng	1486680	4	17.203	3603800	20.0
4H-Cyclopenta[d...]	18.274	2.2	ng	389759	4	17.203	3603800	20.0
Phenol, 2-[(4-h...	18.603	7.0	ng	1252110	4	17.203	3603800	20.0
Phenol, 4,4'-me...	19.015	6.3	ng	1126670	4	17.203	3603800	20.0
Octadecanoic acid	19.374	2.1	ng	697492	5	21.380	6729310	20.0
1,1'-Biphenyl, ...	20.845	2.5	ng	843517	5	21.380	6729310	20.0
(Z)-3-(pentadec...	21.092	2.1	ng	708134	5	21.380	6729310	20.0
1,1'-Biphenyl, ...	21.774	3.5	ng	1177520	5	21.380	6729310	20.0
2,2',3,4,4',5,5...	21.844	3.5	ng	1165390	5	21.380	6729310	20.0