

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
 Data File : BM037439.D
 Acq On : 09 Nov 2022 16:29
 Operator : CG/JU
 Sample : N5436-13 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 250FERRYBLVD-TFS-13

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: BM037439.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	321	rBV	165451	226872	1.70%	0.317%
2	5.398	386	393	398	rBV	527463	755380	5.67%	1.056%
3	7.004	661	666	677	rVB	458920	784307	5.89%	1.097%
4	7.822	799	805	814	rVB	1103627	1670210	12.54%	2.335%
5	8.610	928	939	956	rBV	136305	337143	2.53%	0.471%
6	8.992	997	1004	1009	rBV	295227	498187	3.74%	0.697%
7	9.810	1134	1143	1154	rBV	243764	486737	3.65%	0.681%
8	10.063	1174	1186	1193	rBV2	93373	264586	1.99%	0.370%
9	10.622	1270	1281	1286	rBV	1413021	2320394	17.42%	3.245%
10	10.675	1286	1290	1297	rVV	187750	308050	2.31%	0.431%
11	11.457	1413	1423	1428	rBV2	66139	138817	1.04%	0.194%
12	12.286	1558	1564	1570	rBV	88634	142922	1.07%	0.200%
13	12.980	1677	1682	1691	rBV	102896	177526	1.33%	0.248%
14	13.086	1694	1700	1711	rVB	790374	1154460	8.67%	1.614%
15	13.298	1728	1736	1745	rVB4	65536	148531	1.11%	0.208%
16	13.945	1837	1846	1856	rBV4	128716	263694	1.98%	0.369%
17	14.463	1924	1934	1940	rBV	2239134	3191816	23.96%	4.463%
18	14.527	1940	1945	1953	rVB	216447	355190	2.67%	0.497%
19	14.862	1996	2002	2011	rBV	120168	232419	1.74%	0.325%
20	15.509	2106	2112	2122	rBV	190836	299019	2.24%	0.418%
21	15.709	2142	2146	2151	rBV4	93717	139566	1.05%	0.195%
22	15.951	2182	2187	2193	rBV	676754	951017	7.14%	1.330%
23	17.204	2395	2400	2404	rBV	2576253	3619556	27.17%	5.061%
24	17.245	2404	2407	2414	rVB	668083	963075	7.23%	1.347%
25	17.339	2419	2423	2427	rVB	146458	190745	1.43%	0.267%
26	17.851	2505	2510	2517	rVB	2223736	2710240	20.34%	3.790%
27	18.098	2546	2552	2563	rBV2	957971	1873272	14.06%	2.619%
28	18.274	2578	2582	2587	rBV2	199080	332547	2.50%	0.465%
29	18.392	2598	2602	2606	rBV2	227305	329562	2.47%	0.461%
30	18.603	2633	2638	2643	rBV	763687	1098643	8.25%	1.536%
31	18.980	2699	2702	2705	rBV2	241773	349827	2.63%	0.489%
32	19.015	2705	2708	2718	rVB	829585	1227526	9.21%	1.716%
33	19.262	2746	2750	2755	rBV	1341308	1825324	13.70%	2.552%
34	19.315	2755	2759	2764	rVB	494040	654567	4.91%	0.915%
35	19.380	2766	2770	2773	rBV	704844	915304	6.87%	1.280%
36	19.627	2804	2812	2820	rVB	11330923	13323160	100.00%	18.630%

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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	19.827	2842	2846	2849	rBV	1352587	1580243	11.86%	2.210%
38	19.868	2849	2853	2859	rVB	5484643	6735201	50.55%	9.418%
39	20.268	2918	2921	2927	rBV2	412192	503536	3.78%	0.704%
40	20.550	2966	2969	2972	rBV	367206	462099	3.47%	0.646%
41	20.586	2973	2975	2979	rVB	375865	400281	3.00%	0.560%
42	20.844	3016	3019	3025	rVB2	766219	914903	6.87%	1.279%
43	21.003	3044	3046	3049	rBV	438631	476653	3.58%	0.667%
44	21.091	3059	3061	3069	rVB	959025	1668693	12.52%	2.333%
45	21.386	3105	3111	3122	rBV2	3528061	6689489	50.21%	9.354%
46	21.780	3175	3178	3185	rVB2	1104670	1333588	10.01%	1.865%
47	21.850	3187	3190	3193	rBV	972990	1118070	8.39%	1.563%
48	22.909	3367	3370	3375	rVB	774605	1041152	7.81%	1.456%
49	23.727	3504	3509	3523	rVB2	2240975	4331532	32.51%	6.057%

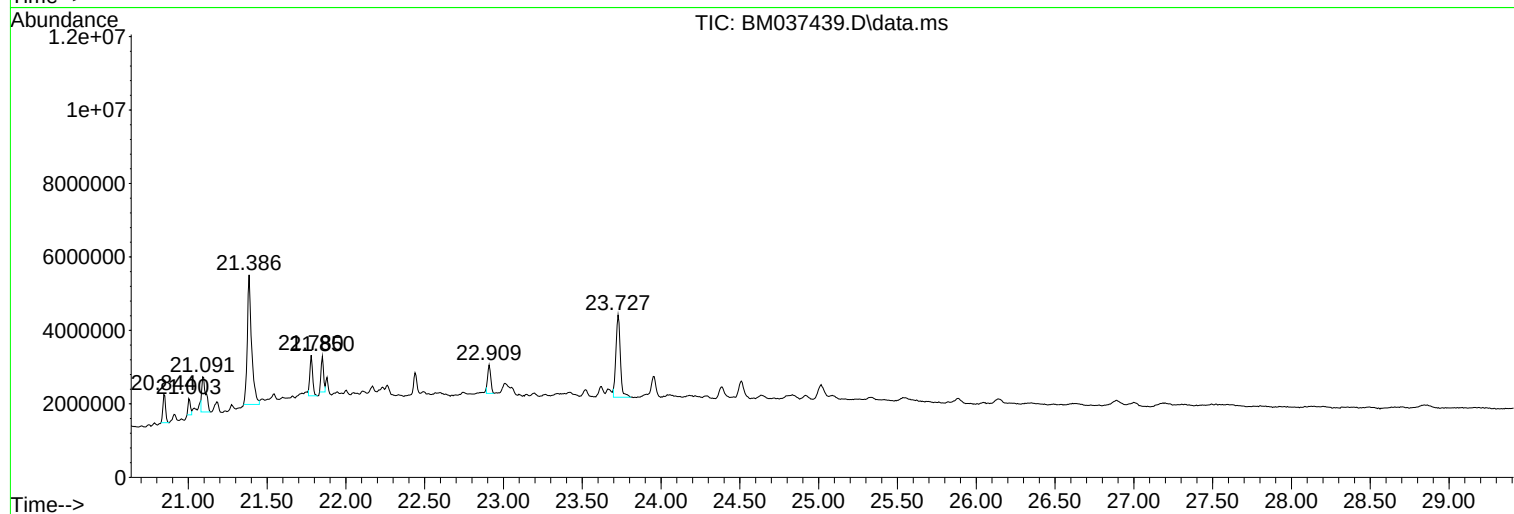
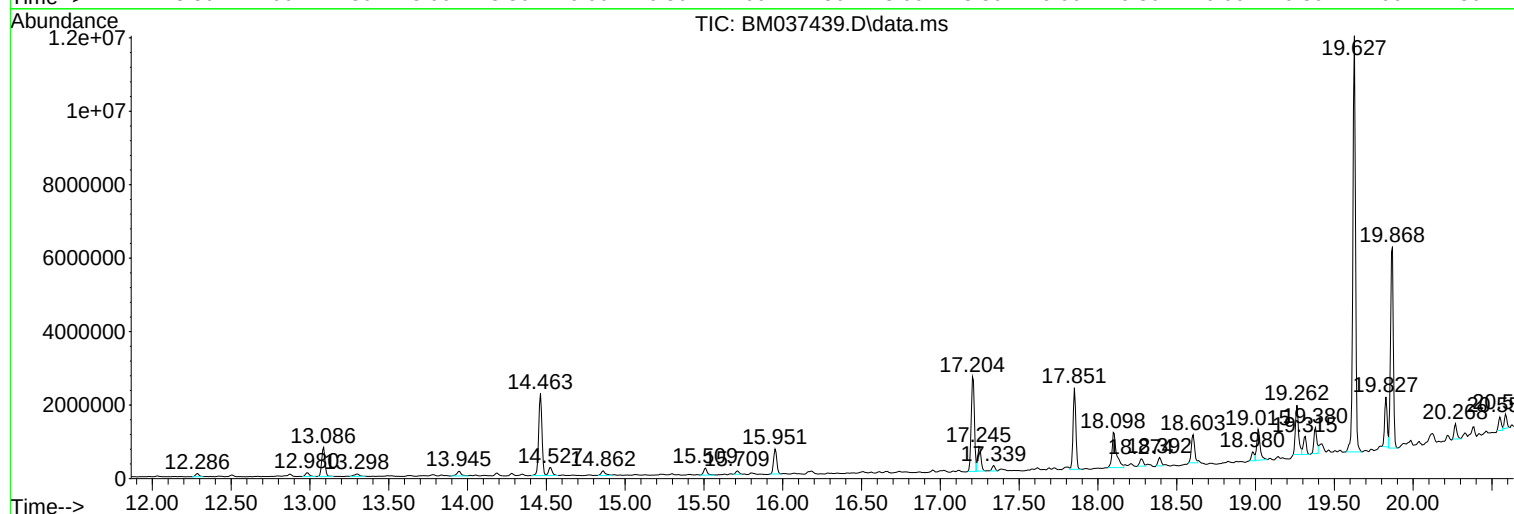
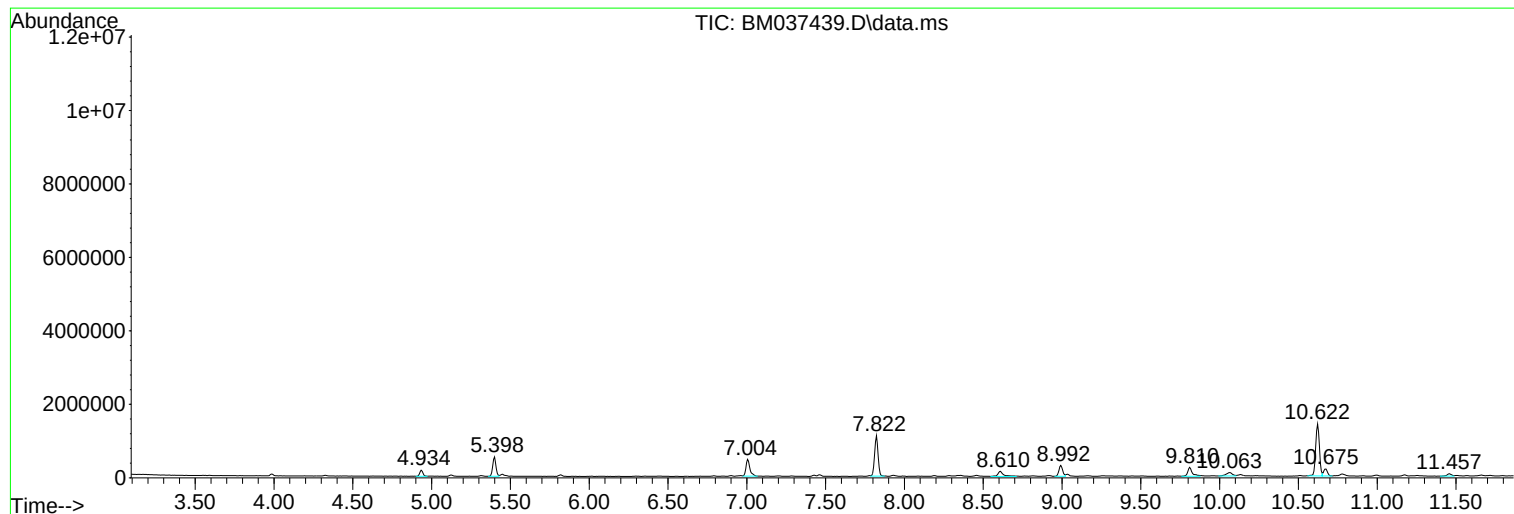
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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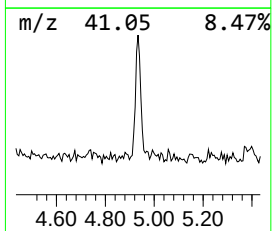
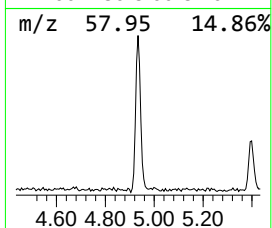
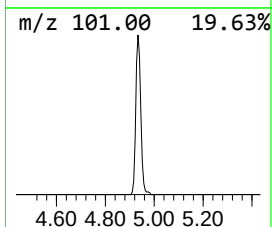
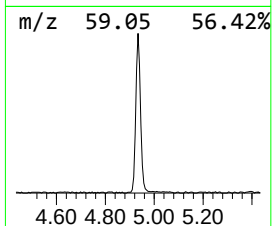
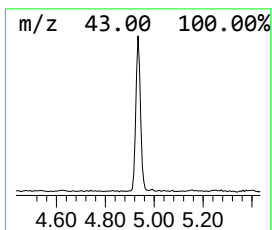
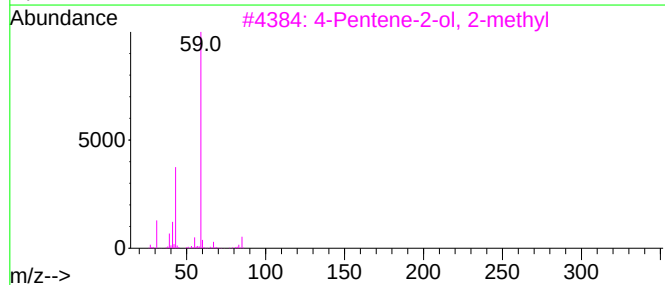
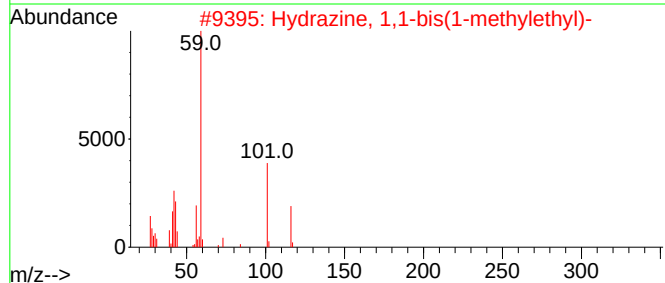
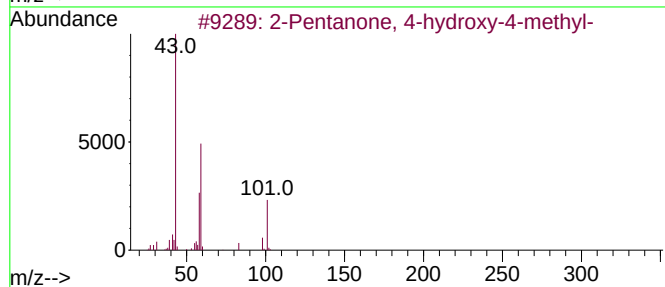
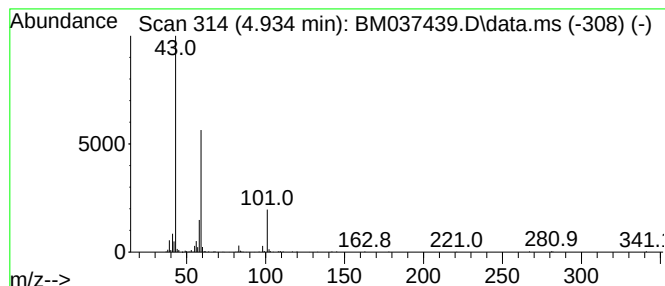
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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	2.72 ng	226872	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	50		
2	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	33		
3	4-Pentene-2-ol, 2-methyl	100	C6H12O	000624-97-5	17		
4	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	16		
5	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	12		



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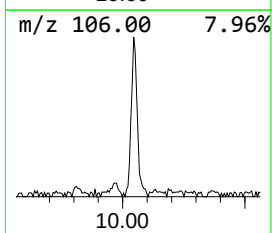
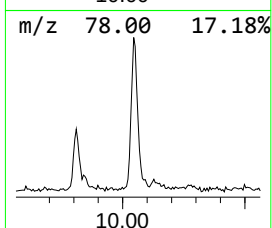
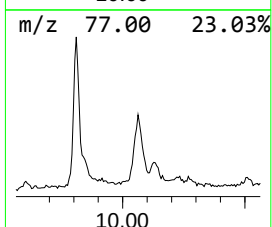
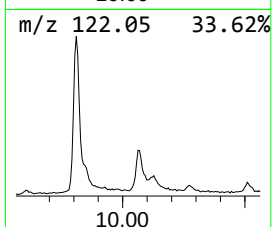
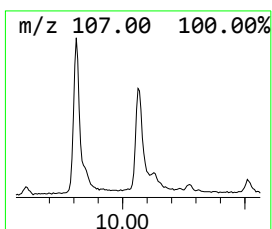
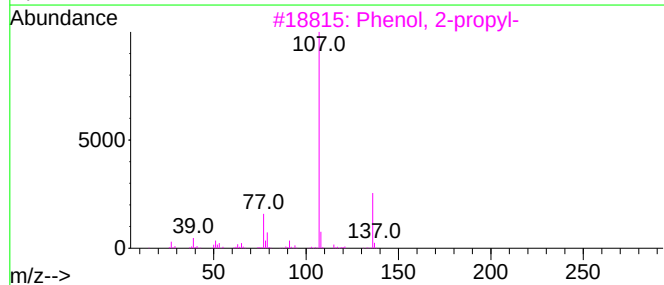
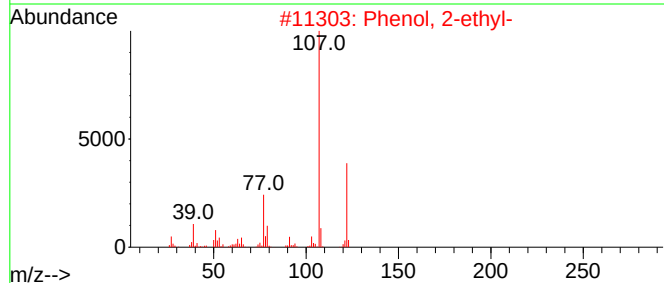
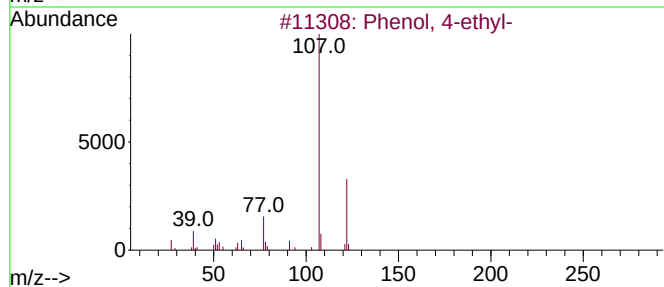
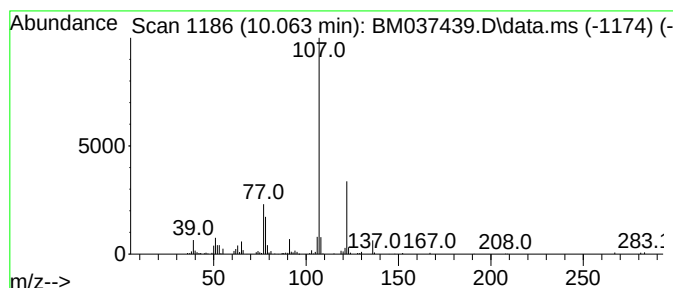
Instrument :
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ClientSampleId :
250FERRYBLVD-TFS-13

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 Phenol, 4-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.063	2.28 ng	264586	Naphthalene-d8	10.622	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-ethyl-	122	C8H10O	000123-07-9	93
2	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	83
3	Phenol, 2-propyl-	136	C9H12O	000644-35-9	76
4	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	72
5	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	72



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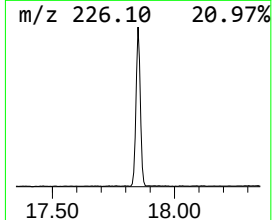
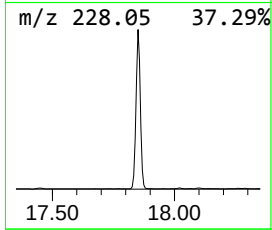
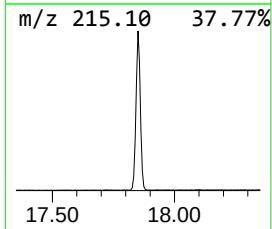
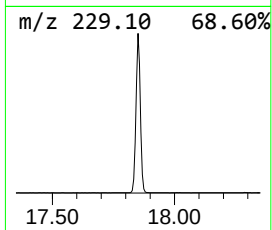
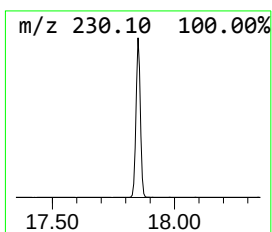
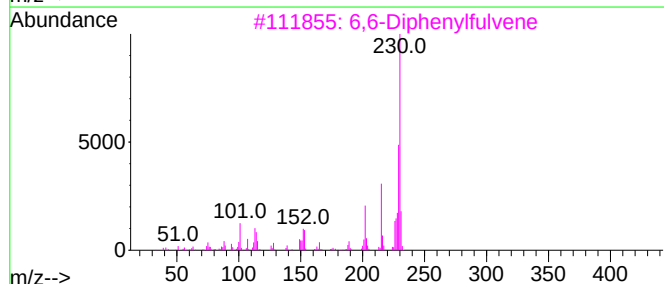
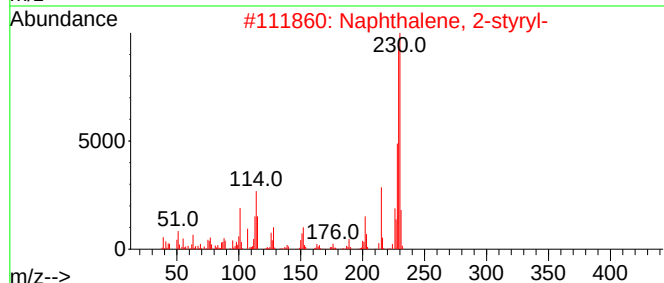
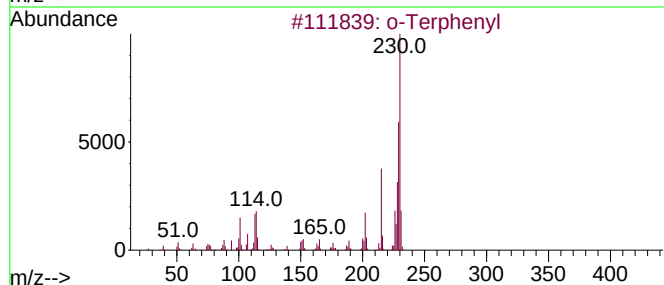
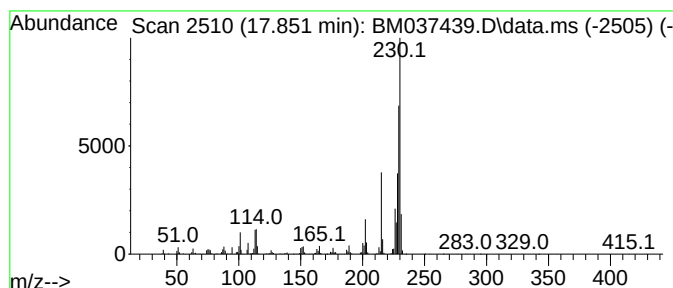
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 o-Terphenyl Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.851	14.98 ng	2710240	Phenanthrene-d10	17.203

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	93
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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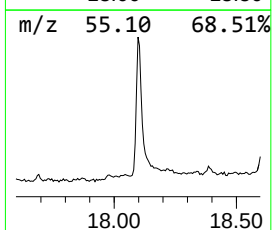
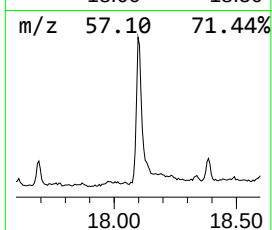
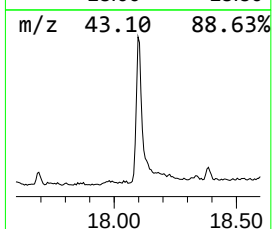
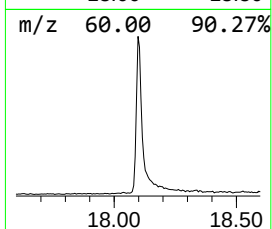
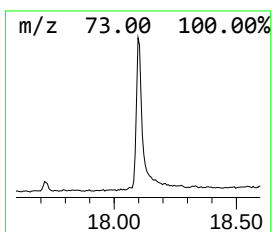
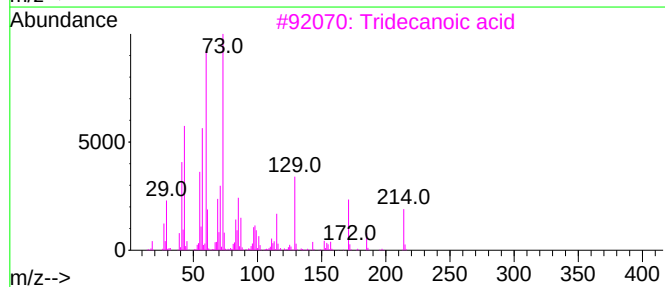
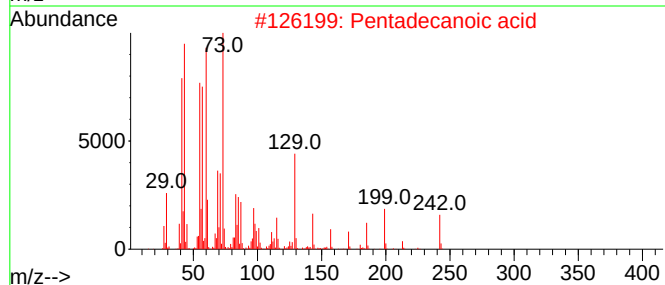
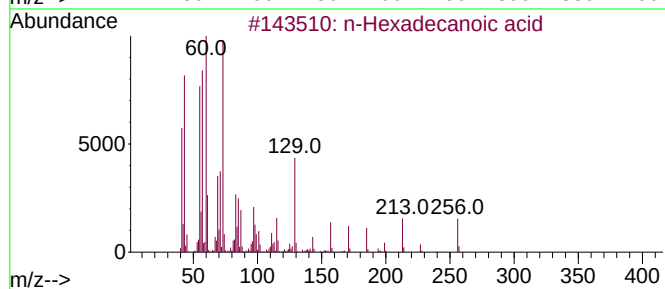
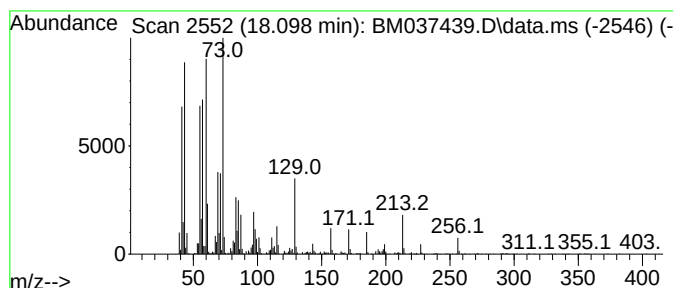
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.098	10.35 ng	1873270	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	90
3	Tridecanoic acid	214	C13H26O2	000638-53-9	81
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5	Octadecanoic acid	284	C18H36O2	000057-11-4	64



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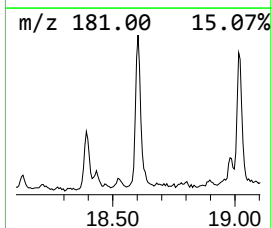
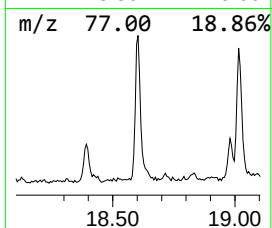
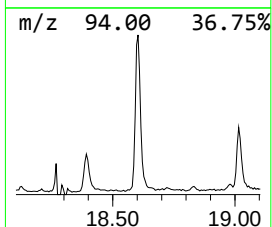
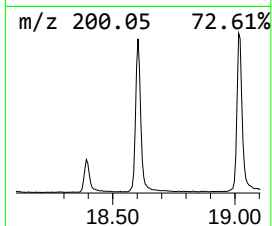
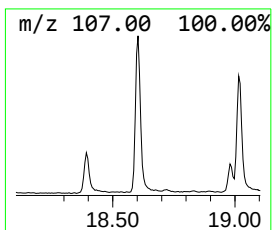
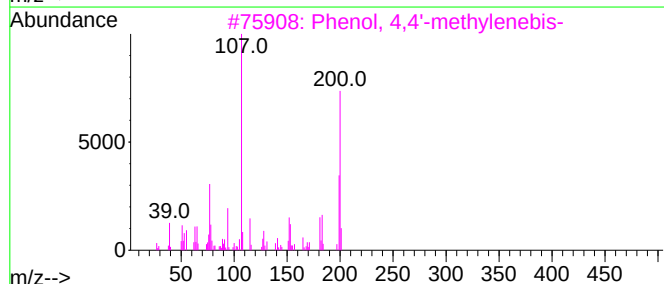
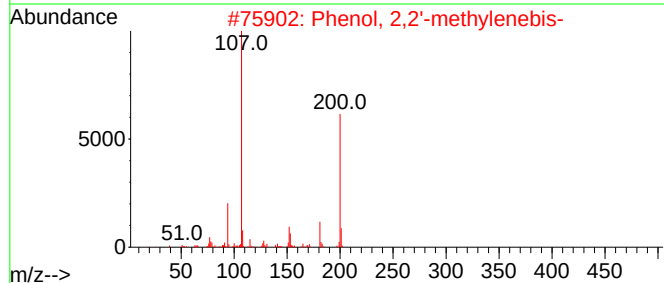
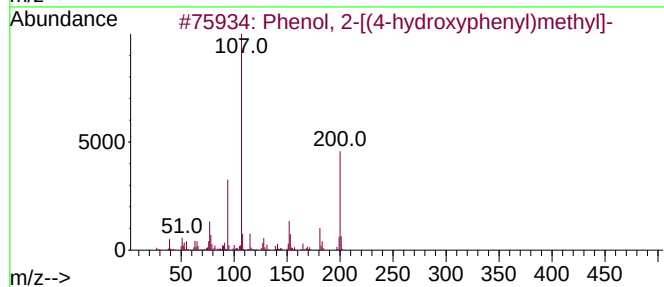
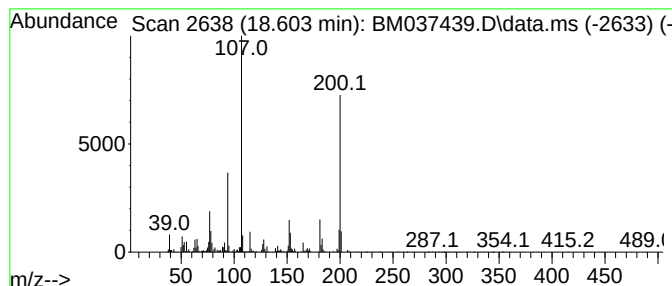
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phenol, 2-[(4-hydroxyphenyl)... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.603	6.07 ng	1098640	Phenanthrene-d10	17.203

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	93
3			Phenol, 4,4'-methylenebis-	200	C13H12O2	000620-92-8	80
4			2,2'-Methylenediphenol, acetate	242	C15H14O3	1000462-45-6	64
5			Diamidafos	200	C8H13N2O2P	001754-58-1	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037439.D
Acq On : 09 Nov 2022 16:29
Operator : CG/JU
Sample : N5436-13 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-13

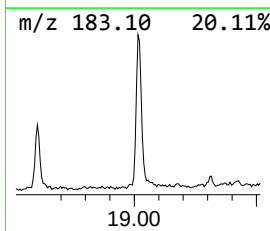
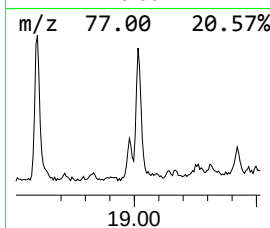
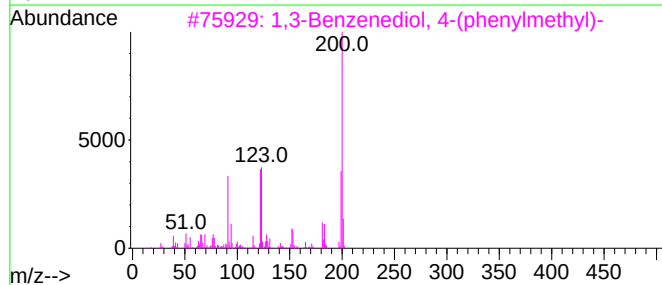
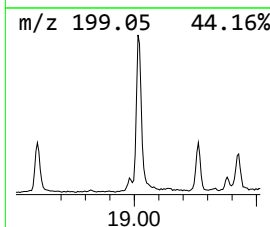
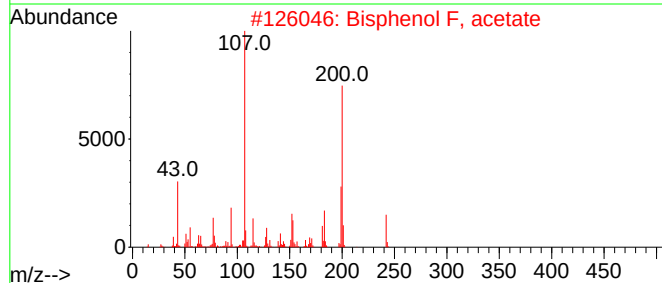
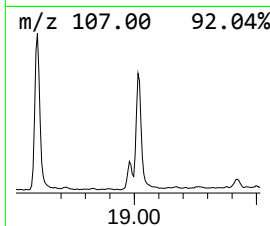
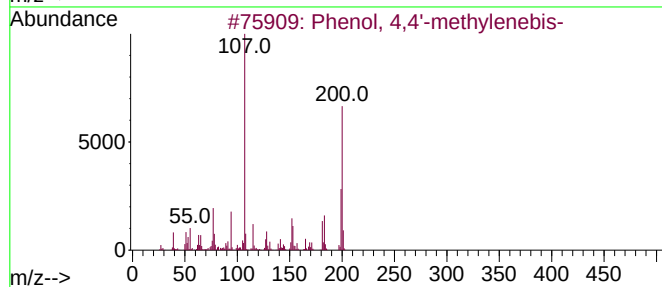
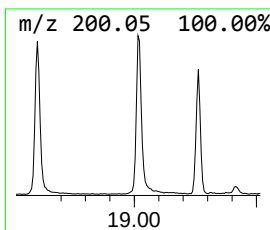
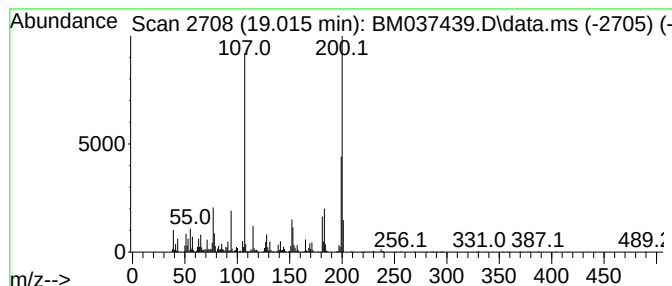
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenol, 4,4'-methylenebis- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.015	6.78 ng	1227530	Phenanthrene-d10	17.203

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4,4'-methylenebis-	200	C13H12O2	000620-92-8	98
2			Bisphenol F, acetate	242	C15H14O3	035250-15-8	87
3			1,3-Benzenediol, 4-(phenylmethyl)-	200	C13H12O2	002284-30-2	83
4			Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	76
5			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037439.D
Acq On : 09 Nov 2022 16:29
Operator : CG/JU
Sample : N5436-13 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
250FERRYBLVD-TFS-13

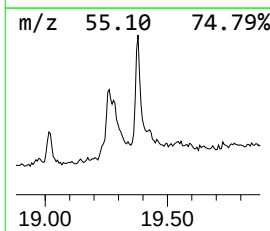
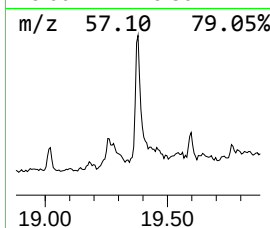
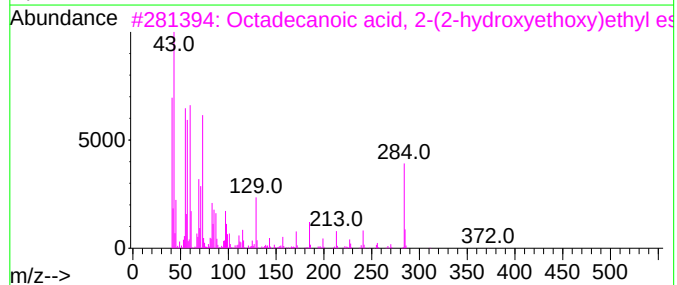
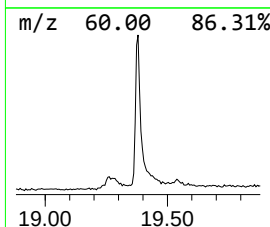
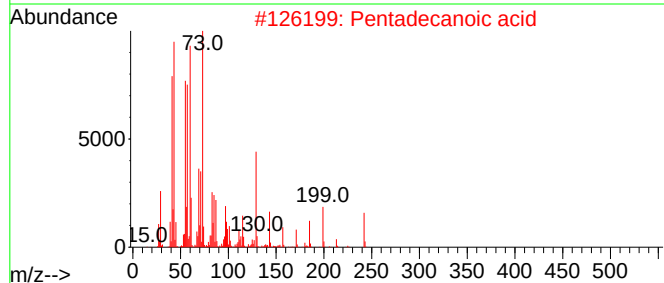
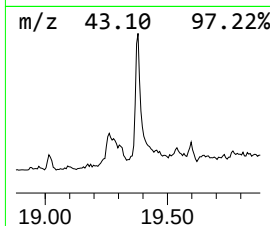
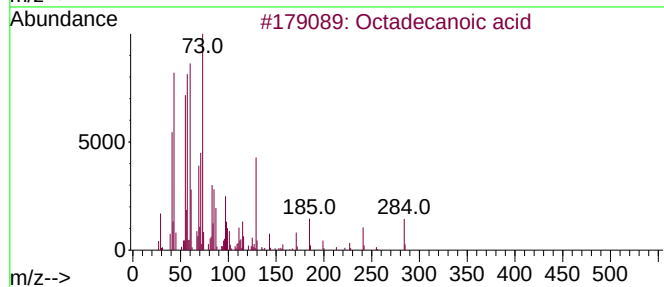
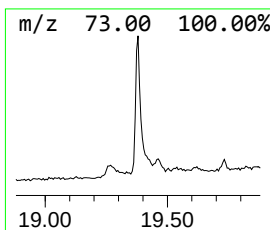
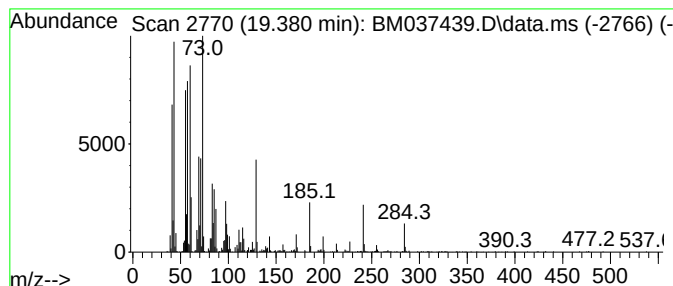
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.380	2.74 ng	915304	Chrysene-d12	21.386

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	93
3			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	90
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	87
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037439.D
Acq On : 09 Nov 2022 16:29
Operator : CG/JU
Sample : N5436-13 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

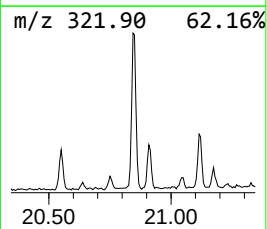
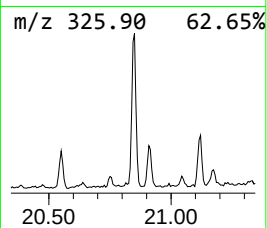
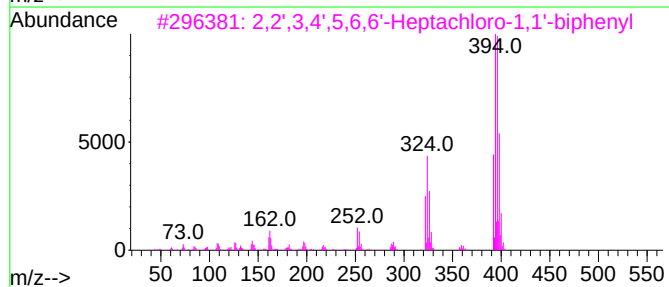
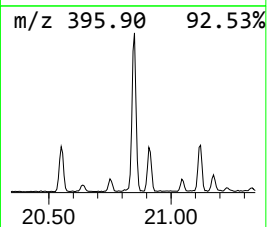
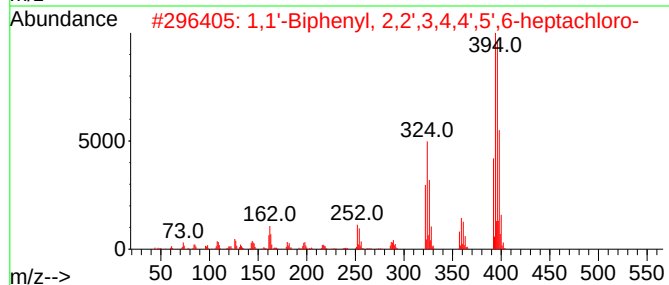
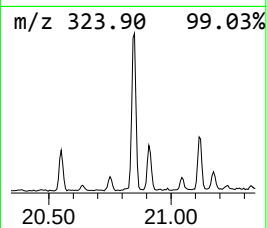
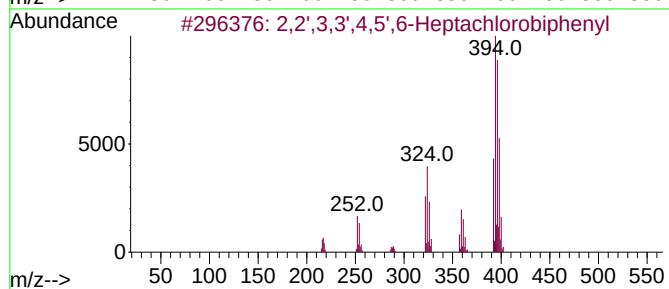
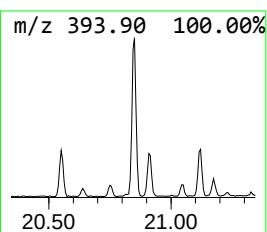
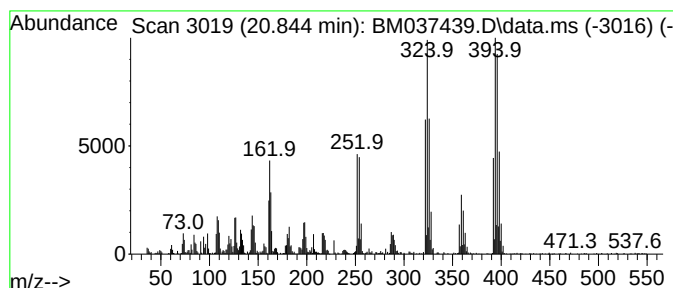
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2,2',3,3',4,5',6-Heptachlor... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.844	2.74 ng	914903	Chrysene-d12	21.386	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,3',4,5',6-Heptachlorobiph...	392	C12H3Cl7	040186-70-7	99
2	1,1'-Biphenyl, 2,2',3,4,4',5',6-...	392	C12H3Cl7	052663-69-1	99
3	2,2',3,4',5,6,6'-Heptachloro-1,1...	392	C12H3Cl7	074487-85-7	99
4	1,1'-Biphenyl, 2,3,3',4,4',5,5'-...	392	C12H3Cl7	039635-31-9	99
5	1,1'-Biphenyl, 2,3,3',4',5,5',6-...	392	C12H3Cl7	069782-91-8	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037439.D
Acq On : 09 Nov 2022 16:29
Operator : CG/JU
Sample : N5436-13 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
250FERRYBLVD-TFS-13

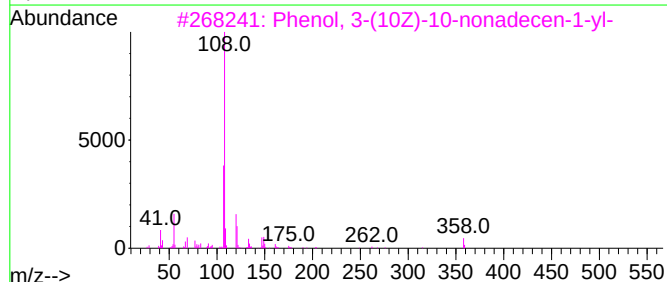
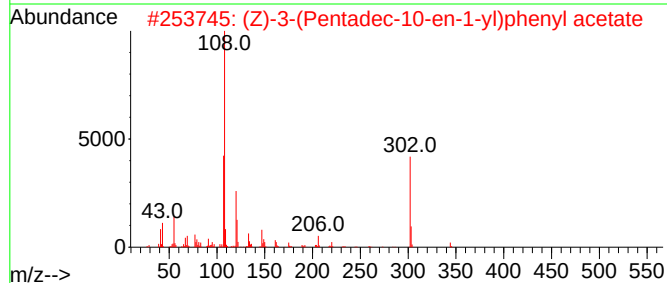
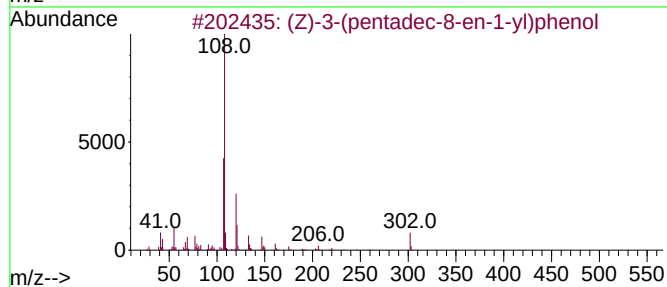
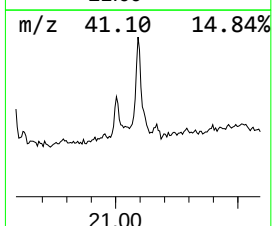
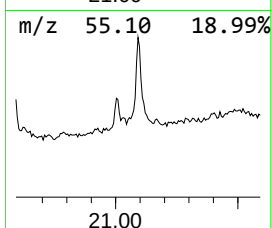
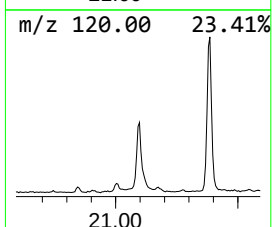
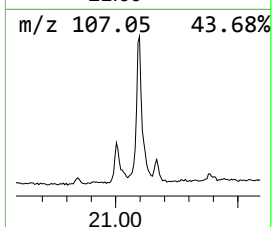
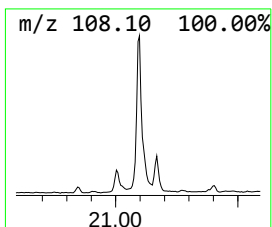
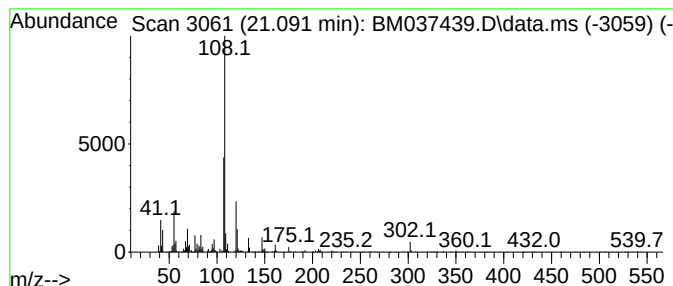
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 (Z)-3-(pentadec-8-en-1-yl)p... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.092	4.99 ng	1668690	Chrysene-d12	21.386

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		(Z)-3-(Pentadec-10-en-1-yl)pheny...	344	C23H36O2	1000478-00-3	94
3		Phenol, 3-(10Z)-10-nonadecen-1-yl-	358	C25H42O	936109-24-9	72
4		Phenol, 3-undecyl-	248	C17H28O	020056-72-8	64
5		Phenol, 3-pentadecyl-	304	C21H36O	000501-24-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037439.D
Acq On : 09 Nov 2022 16:29
Operator : CG/JU
Sample : N5436-13 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

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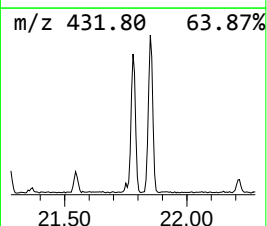
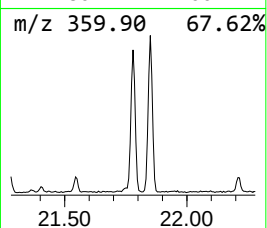
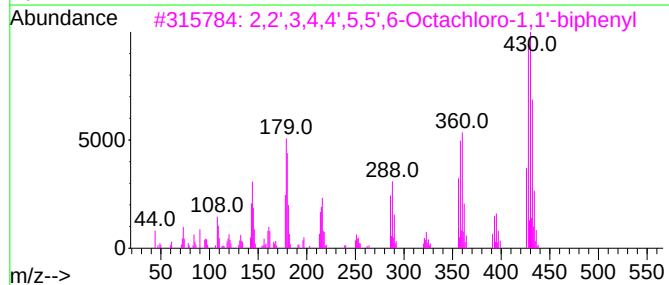
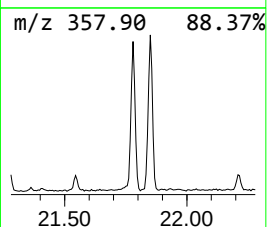
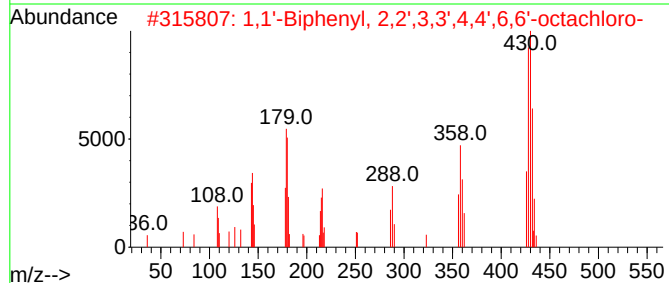
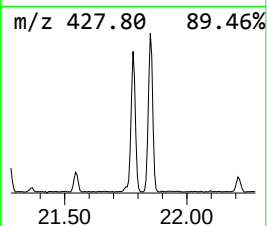
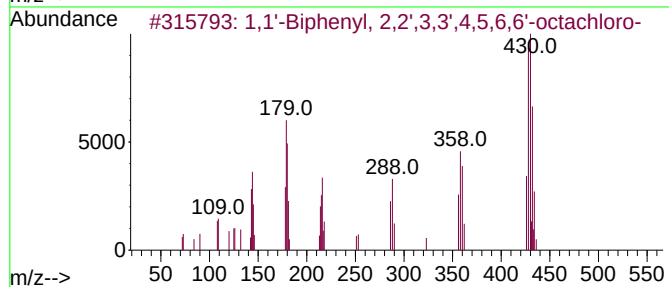
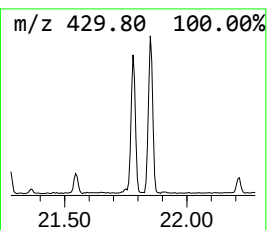
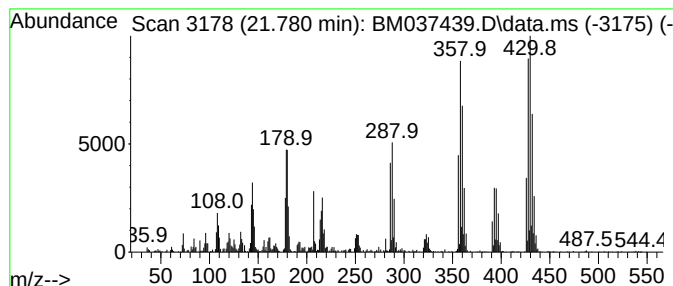
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.780	3.99 ng	1333590	Chrysene-d12	21.386

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	2,2',3,4,4',5,5',6-Octachloro-1...	426	C12H2Cl8	052663-76-0	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',5...	426	C12H2Cl8	042740-50-1	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
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Misc :
ALS Vial : 8 Sample Multiplier: 1

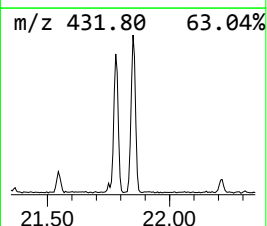
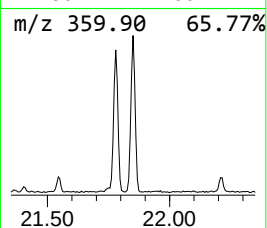
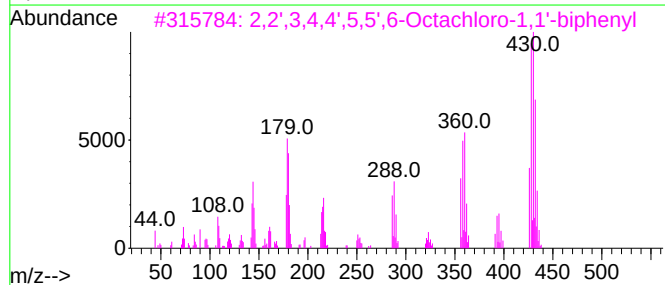
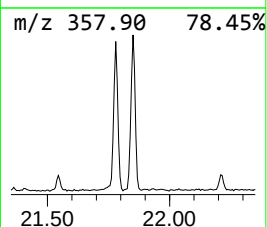
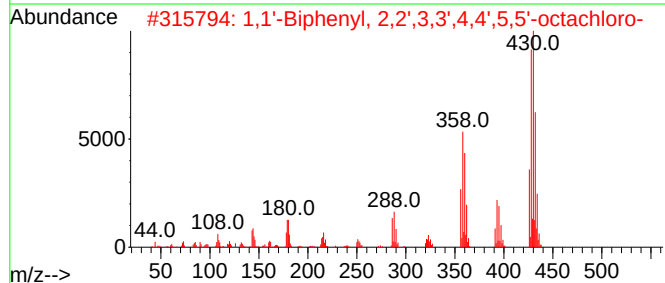
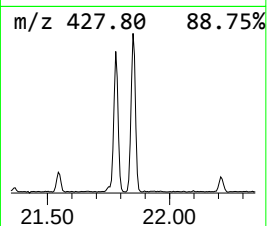
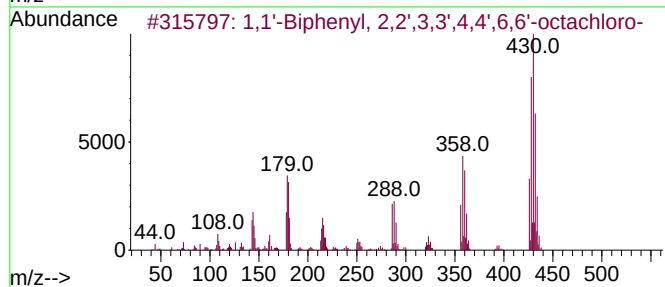
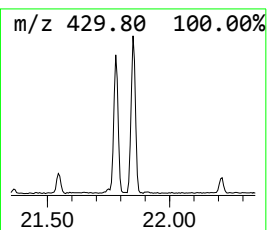
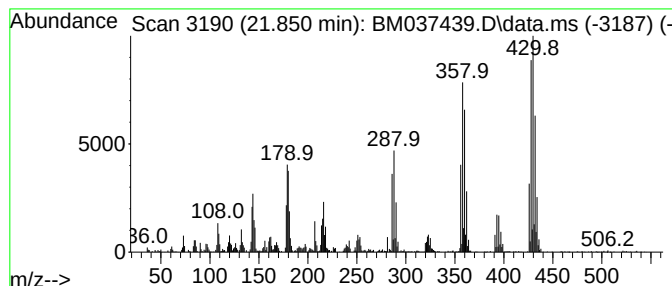
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.850	3.34 ng	1118070	Chrysene-d12	21.386	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,3',4,4',6,...	426	C12H2Cl8	033091-17-7	99
2	1,1'-Biphenyl, 2,2',3,3',4,4',5,...	426	C12H2Cl8	035694-08-7	99
3	2,2',3,4,4',5,5',6-Octachloro-1,...	426	C12H2Cl8	052663-76-0	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',4,5,6,6...	426	C12H2Cl8	052663-73-7	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037439.D
Acq On : 09 Nov 2022 16:29
Operator : CG/JU
Sample : N5436-13 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-13

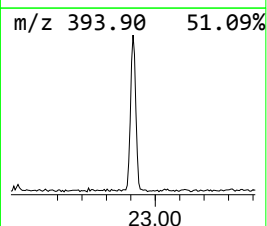
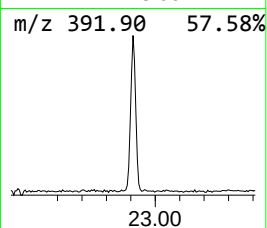
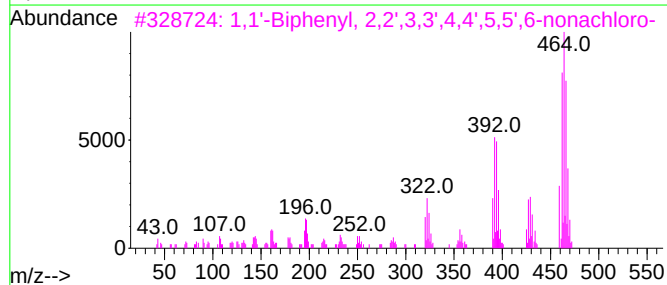
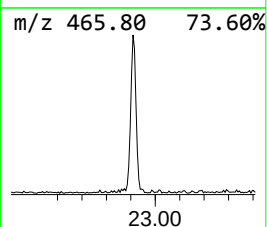
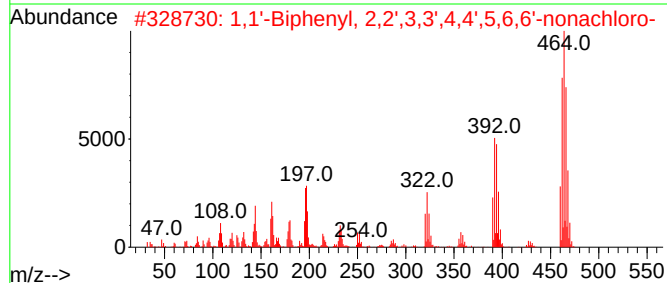
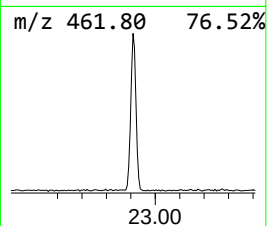
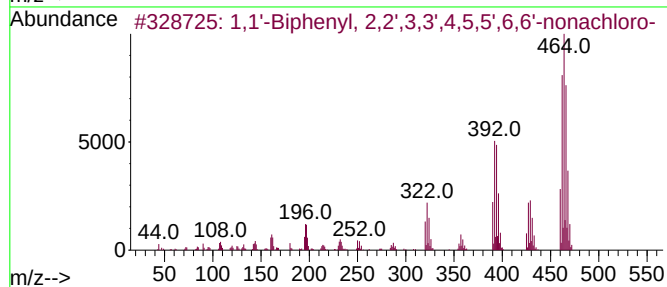
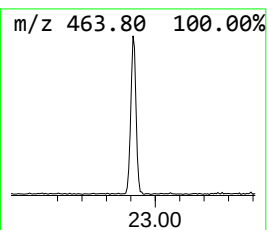
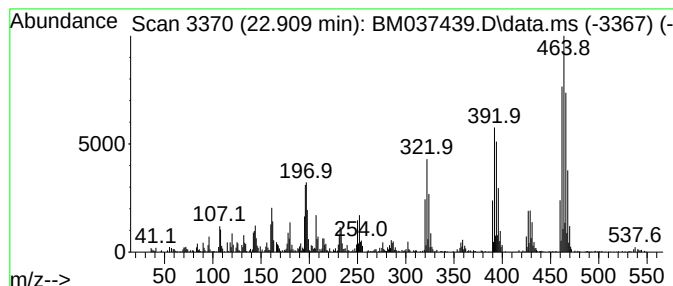
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.909	4.81 ng	1041150	Perylene-d12	23.727

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,3',4,5,5',...	460	C12HC19	052663-77-1	99	
2	1,1'-Biphenyl, 2,2',3,3',4,4',5,...	460	C12HC19	052663-79-3	99	
3	1,1'-Biphenyl, 2,2',3,3',4,4',5,...	460	C12HC19	040186-72-9	95	
4	Spiro[benzofuran-2(3H),1'-[2,5]c...	462	C14H9Br3O3	052498-93-8	35	
5	Spiro[9,9']difluorene, 2,2'-(1,4...	462	C31H26O4	1000191-82-0	9	



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037439.D
Acq On : 09 Nov 2022 16:29
Operator : CG/JU
Sample : N5436-13 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-13

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	2.7	ng	226872	1	7.822	1670210	20.0
Phenol, 4-ethyl-	10.063	2.3	ng	264586	2	10.622	2320390	20.0
o-Terphenyl	17.851	15.0	ng	2710240	4	17.203	3619560	20.0
n-Hexadecanoic ...	18.098	10.4	ng	1873270	4	17.203	3619560	20.0
Phenol, 2-[(4-h...	18.603	6.1	ng	1098640	4	17.203	3619560	20.0
Phenol, 4,4'-me...	19.015	6.8	ng	1227530	4	17.203	3619560	20.0
Octadecanoic acid	19.380	2.7	ng	915304	5	21.386	6689490	20.0
2,2',3,3',4,5',...	20.844	2.7	ng	914903	5	21.386	6689490	20.0
(Z)-3-(pentadec...	21.092	5.0	ng	1668690	5	21.386	6689490	20.0
1,1'-Biphenyl, ...	21.780	4.0	ng	1333590	5	21.386	6689490	20.0
1,1'-Biphenyl, ...	21.850	3.3	ng	1118070	5	21.386	6689490	20.0
1,1'-Biphenyl, ...	22.909	4.8	ng	1041150	6	23.727	4331530	20.0