LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037440.D
Acq On : 09 Nov 2022 17:05
Operator : CG/JU
Sample : N5436-04 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-04

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: BM037440.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	321	rBV	116859	159425	1.85%	0.299%
2	5.398	387	393	398	rBV	329317	485230	5.62%	0.910%
3	7.004	660	666	678	rVB	275768	474792	5.50%	0.890%
4	7.822	799	805	814	rVB	1041654	1577426	18.28%	2.958%
5	8.992	997	1004	1009	rBV	180911	316500	3.67%	0.594%
6	9.810	1136	1143	1154	rBV	100124	226310	2.62%	0.424%
7	10.057	1173	1185	1194	rBV5	37976	132812	1.54%	0.249%
8	10.622	1270	1281	1287	rBV	1345436	2217361	25.70%	4.159%
9	10.669	1287	1289	1296	rVB	99129	144660	1.68%	0.271%
10	13.086	1693	1700	1706	rBV	493328	717181	8.31%	1.345%
11	13.298	1728	1736	1744	rVB2	48753	104314	1.21%	0.196%
12	13.939	1837	1845	1852	rBV5	37077	92510	1.07%	0.173%
13	14.186	1882	1887	1892	rBV	66568	99567	1.15%	0.187%
14	14.462	1925	1934	1940	rBV	2071162	3004063	34.82%	5.634%
15	14.521	1940	1944	1952	rVB	340356	516585	5.99%	0.969%
16	14.857	1996	2001	2010	rBV	182619	308986	3.58%	0.579%
17	15.509	2106	2112	2123	rVB	356596	528523	6.13%	0.991%
18	15.709	2142	2146	2151	rVB4	79345	105929	1.23%	0.199%
19	15.951	2182	2187	2194	rBV	469321	669820	7.76%	1.256%
20	17.021	2366	2369	2376	rVB	94769	140433	1.63%	0.263%
21	17.203	2394	2400	2404	rBV	2576053	3550725	41.15%	6.659%
22	17.245	2404	2407	2414	rVB	1290222	1721777	19.95%	3.229%
23	17.339	2419	2423	2427	rVV	192571	260211	3.02%	0.488%
24	17.380	2428	2430	2438	rVB4	44050	97017	1.12%	0.182%
25	17.615	2467	2470	2474	rVB	93180	121074	1.40%	0.227%
26	17.850	2505	2510	2517	rVB	1313152	1699562	19.70%	3.187%
27	18.098	2546	2552	2563	rBV2	758877	1483356	17.19%	2.782%
28	18.274	2578	2582	2586	rBV2	209289	322432	3.74%	0.605%
29	18.386	2597	2601	2606	rBV2	169381	272370	3.16%	0.511%
30	18.603	2632	2638	2643	rBV2	506206	762864	8.84%	1.431%
31	18.980	2699	2702	2705	rBV2	152632	242643	2.81%	0.455%
32	19.015	2705	2708	2718	rVB	498947	739594	8.57%	1.387%
33	19.262	2745	2750	2756	rBV	1170112	1697456	19.67%	3.184%
34	19.309	2756	2758	2763	rVB	250386	319275	3.70%	0.599%
35	19.380	2766	2770	2773	rBV	508707	686098	7.95%	1.287%
36	19.627	2807	2812	2821	rVB	6994494	8628511	100.00%	16.183%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037440.D
Acq On : 09 Nov 2022 17:05
Operator : CG/JU
Sample : N5436-04 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-04

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

37	19.827	2842	2846	2849	rBV	850763	1003653	11.63%	1.882%
38	19.862	2849	2852	2858	rVB	3554862	4292672	49.75%	8.051%
39	20.268	2918	2921	2926	rBV2	222136	305544	3.54%	0.573%
40	20.586	2973	2975	2979	rVB	261455	256934	2.98%	0.482%
41	20.844	3016	3019	3025	rBV3	539447	634985	7.36%	1.191%
42	21.003	3044	3046	3049	rBV	265909	305617	3.54%	0.573%
43	21.091	3059	3061	3071	rVB3	587188	1043544	12.09%	1.957%
44	21.386	3106	3111	3121	rVB	3241985	5459273	63.27%	10.239%
45	21.780	3175	3178	3184	rVB	770407	932310	10.80%	1.749%
46	21.850	3187	3190	3193	rBV	665789	760380	8.81%	1.426%
47	23.727	3504	3509	3516	rVB2	1985361	3697672	42.85%	6.935%

Sum of corrected areas: 53319976

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
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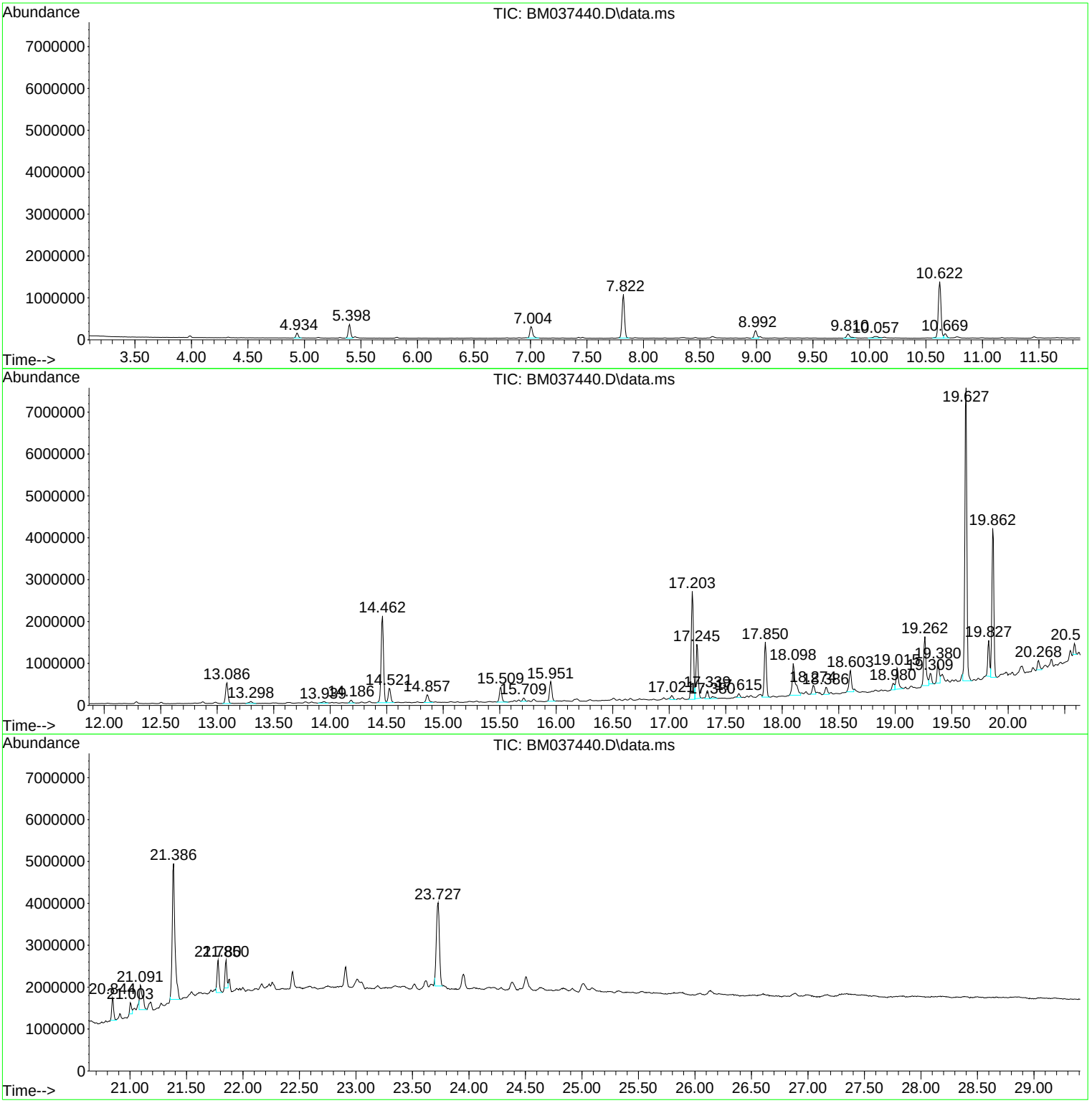
ClientSampleId :
250FERRYBLVD-TFS-04

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037440.D
Acq On : 09 Nov 2022 17:05
Operator : CG/JU
Sample : N5436-04 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-04

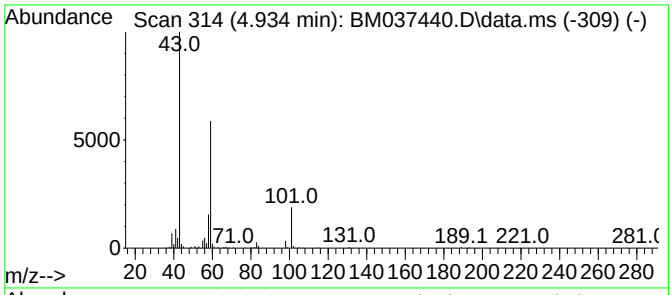
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

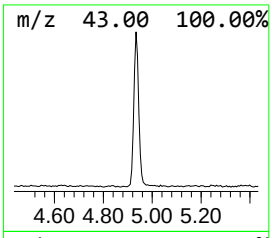
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

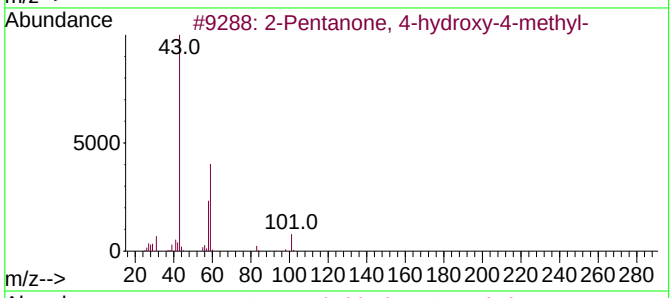
****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

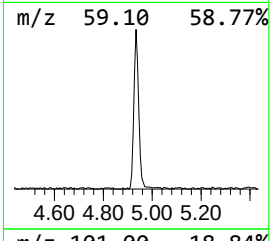
R.T.	EstConc	Area	Relative to ISTD	R.T.
4.934	2.02 ng	159425	1,4-Dichlorobenzene-d4	7.822

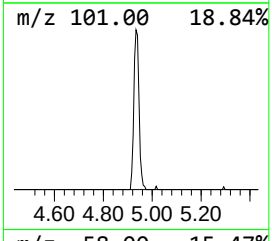
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	32
3	Tetraglyme	222	C10H22O5	000143-24-8	25
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5	Diacetamide	101	C4H7NO2	000625-77-4	9

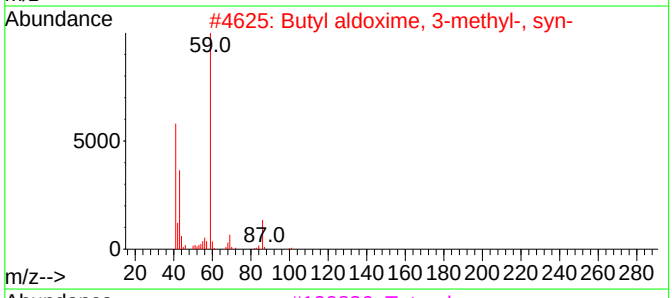


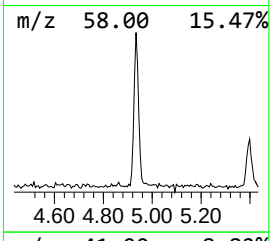


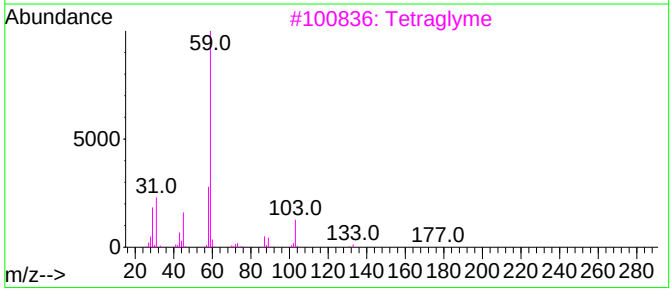


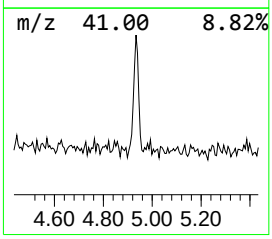












Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037440.D
Acq On : 09 Nov 2022 17:05
Operator : CG/JU
Sample : N5436-04 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-04

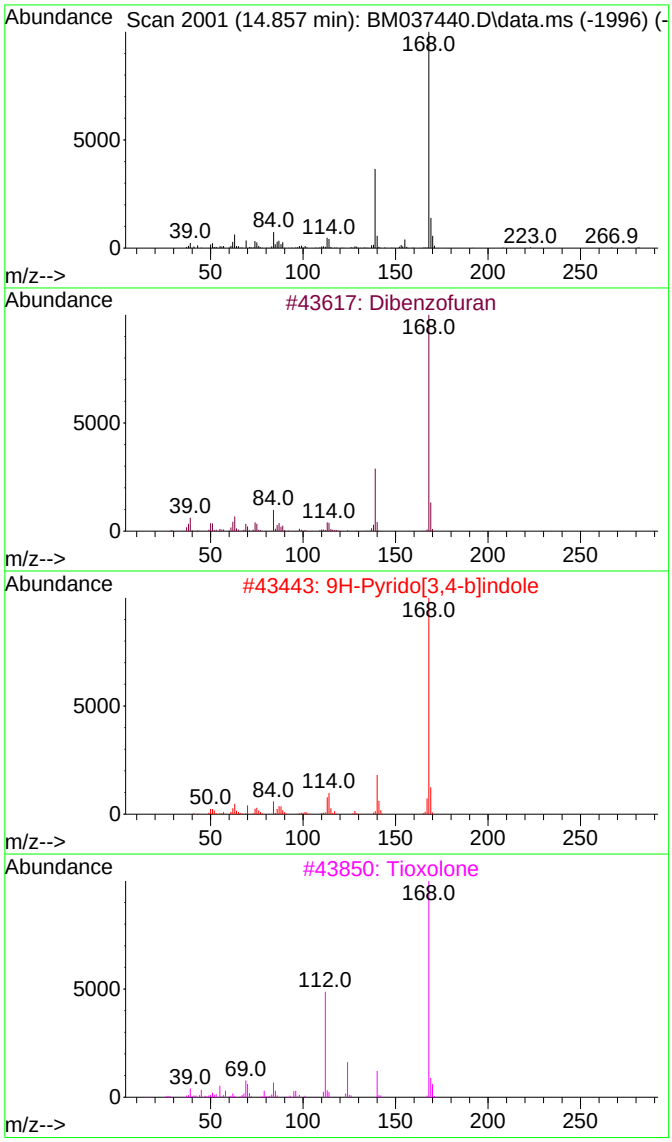
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

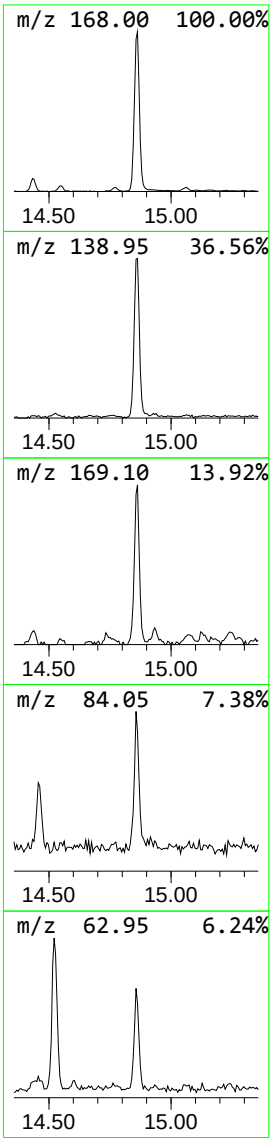
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 9H-Pyrido[3,4-b]indole Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.857	2.06 ng	308986	Acenaphthene-d10	14.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran	168	C12H8O	000132-64-9	90
2	9H-Pyrido[3,4-b]indole	168	C11H8N2	000244-63-3	59
3	Tioxolone	168	C7H4O3S	004991-65-5	59
4	2-Amino-6-fluorobenzothiazole	168	C7H5FN2S	000348-40-3	59
5	1H-Pyrido[2,3-b]indole	168	C11H8N2	000244-76-8	59





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Data File : BM037440.D
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Misc :
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Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-04

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

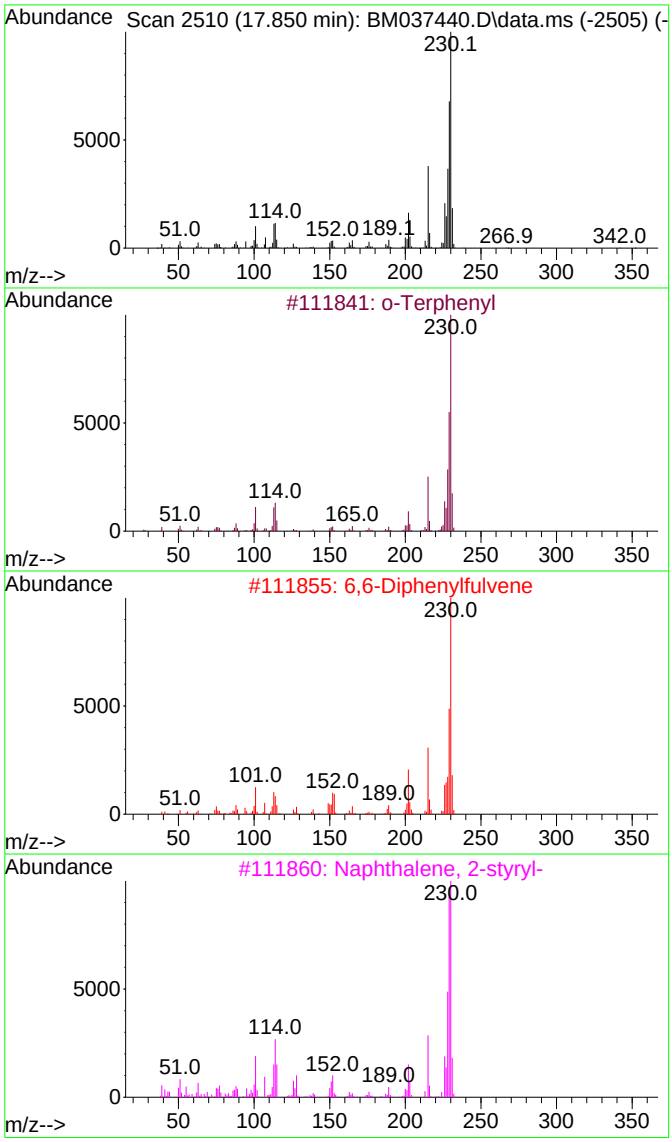
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

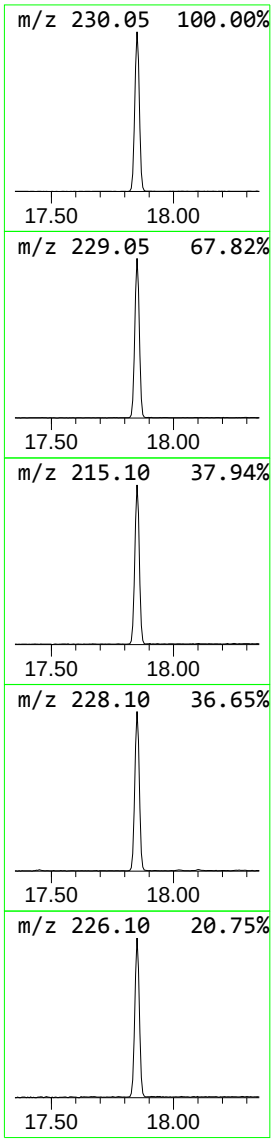
****

Peak Number 3 o-Terphenyl Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.851	9.57 ng	1699560	Phenanthrene-d10	17.203

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Terphenyl	230	C18H14	000084-15-1	98
2	6,6-Diphenylfulvene	230	C18H14	002175-90-8	93
3	Naphthalene, 2-styryl-	230	C18H14	002039-70-5	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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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-04

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

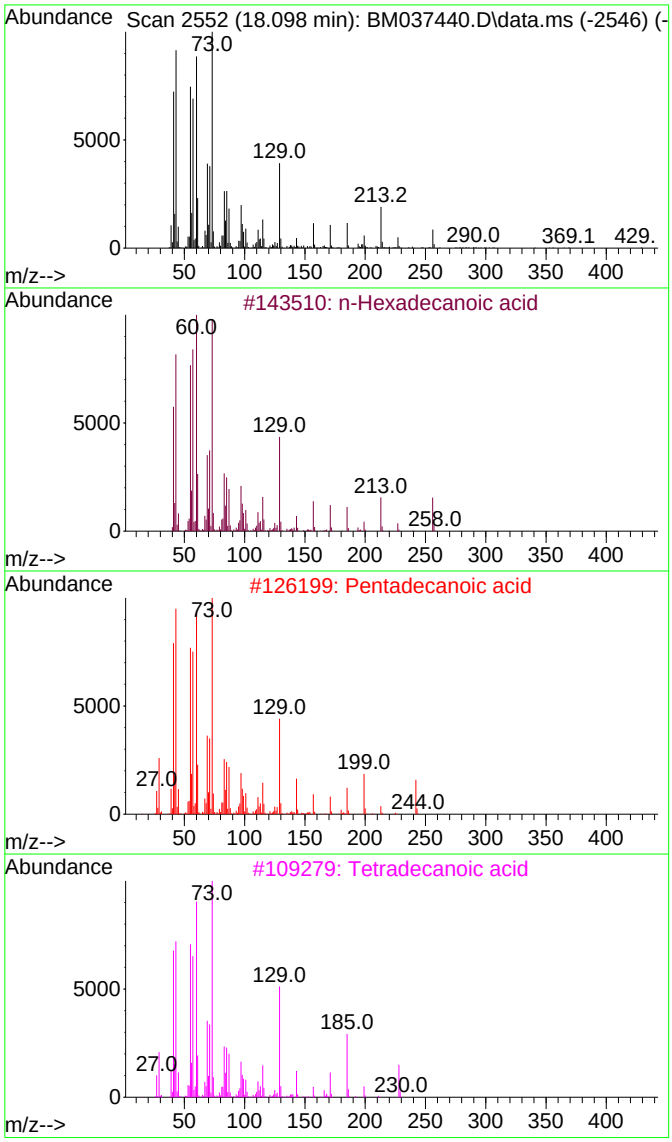
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

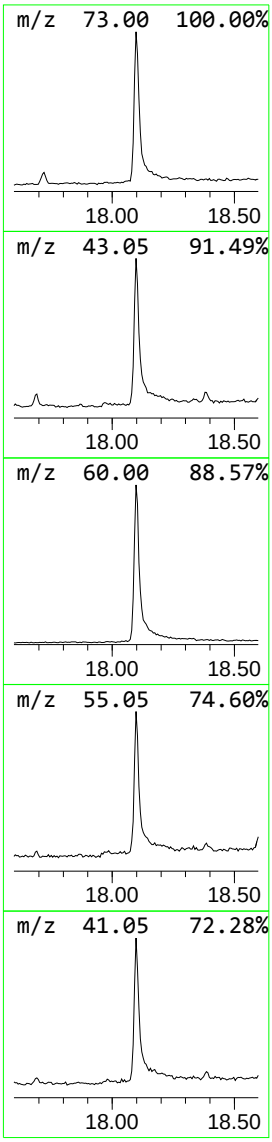
****

Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.098	8.36 ng	1483360	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	87
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4	Octadecanoic acid	284	C18H36O2	000057-11-4	74
5	Tridecanoic acid	214	C13H26O2	000638-53-9	72





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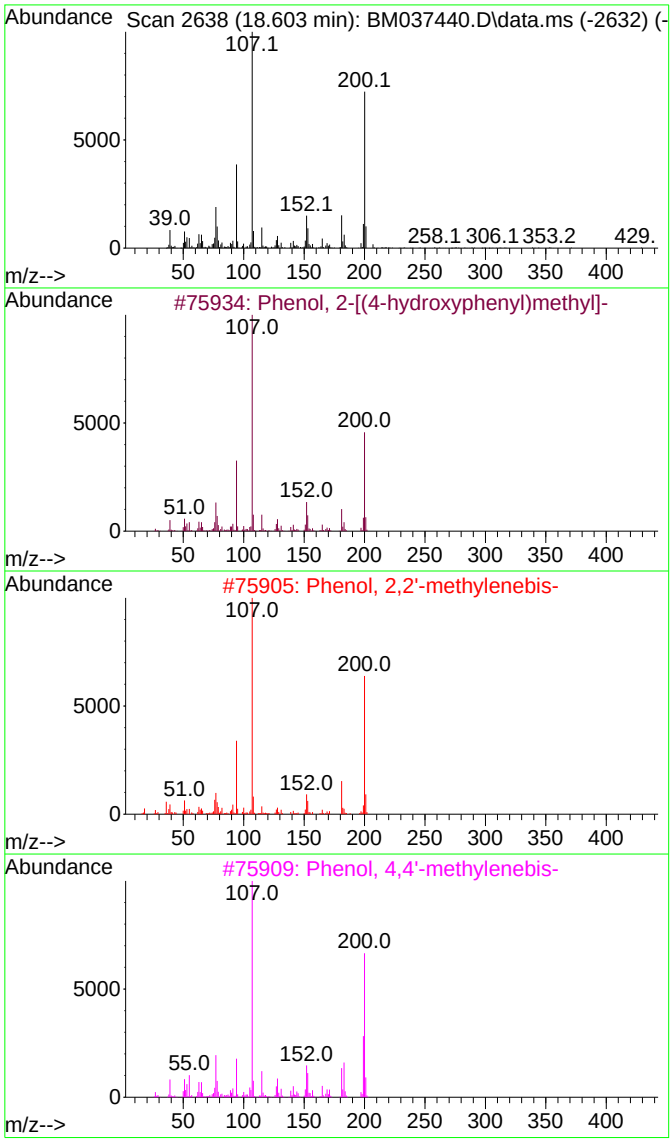
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

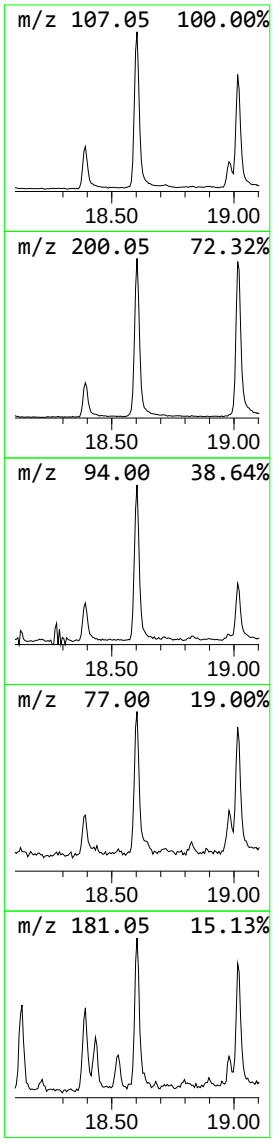
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.603	4.30 ng	762864	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	76
4	2,4'-Dihydroxydiphenylmethane, d...	284	C17H16O4	959033-36-4	72
5	2,2'-Methylenediphenol, diacetate	284	C17H16O4	038782-65-9	64





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

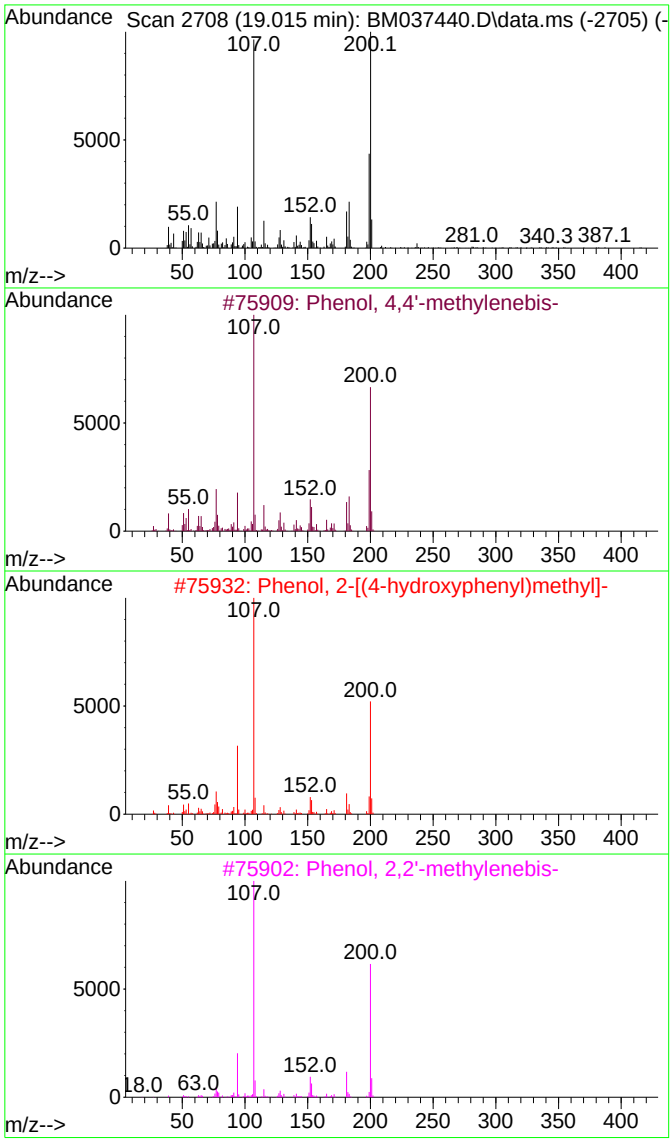
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

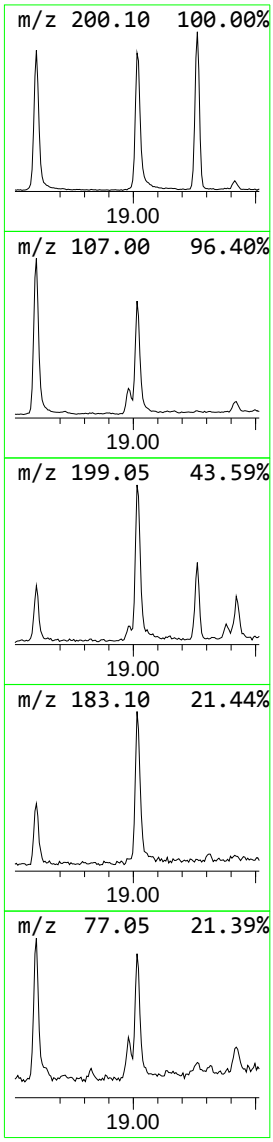
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.015	4.17 ng	739594	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	Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	90
3	Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	81
4	Bisphenol F, acetate	242	C15H14O3	035250-15-8	68
5	Bisphenol F, 2-methylpropionate	270	C17H18O3	1000462-19-4	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
Data File : BM037440.D
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Sample : N5436-04 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
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

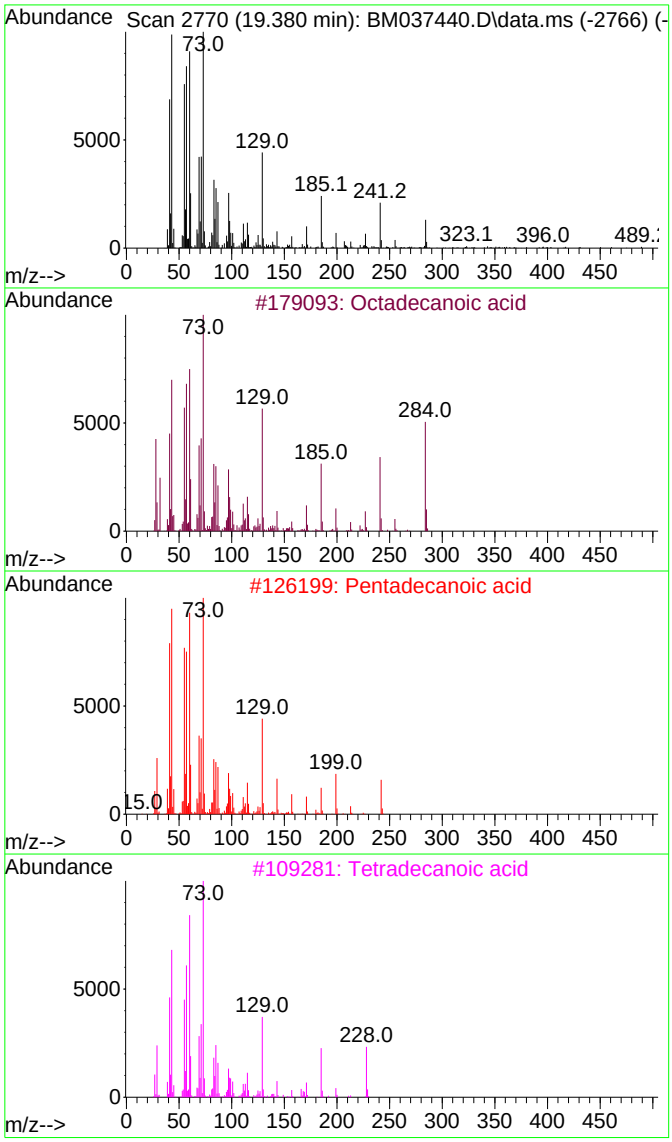
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

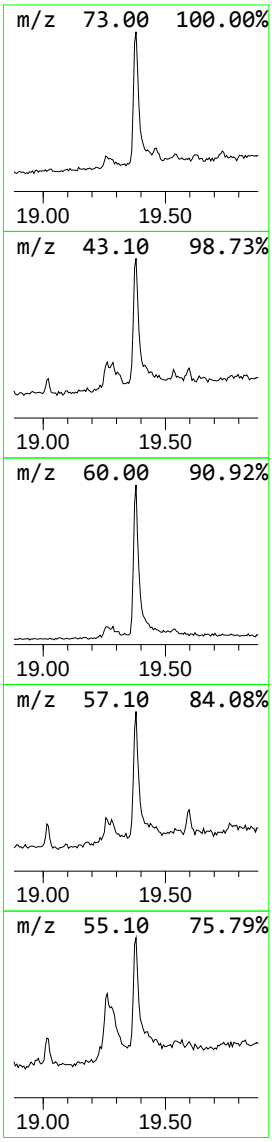
****

Peak Number 7 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.380	2.51 ng	686098	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	91
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
5	Tridecanoic acid	214	C13H26O2	000638-53-9	72





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Data File : BM037440.D
Acq On : 09 Nov 2022 17:05
Operator : CG/JU
Sample : N5436-04 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-04

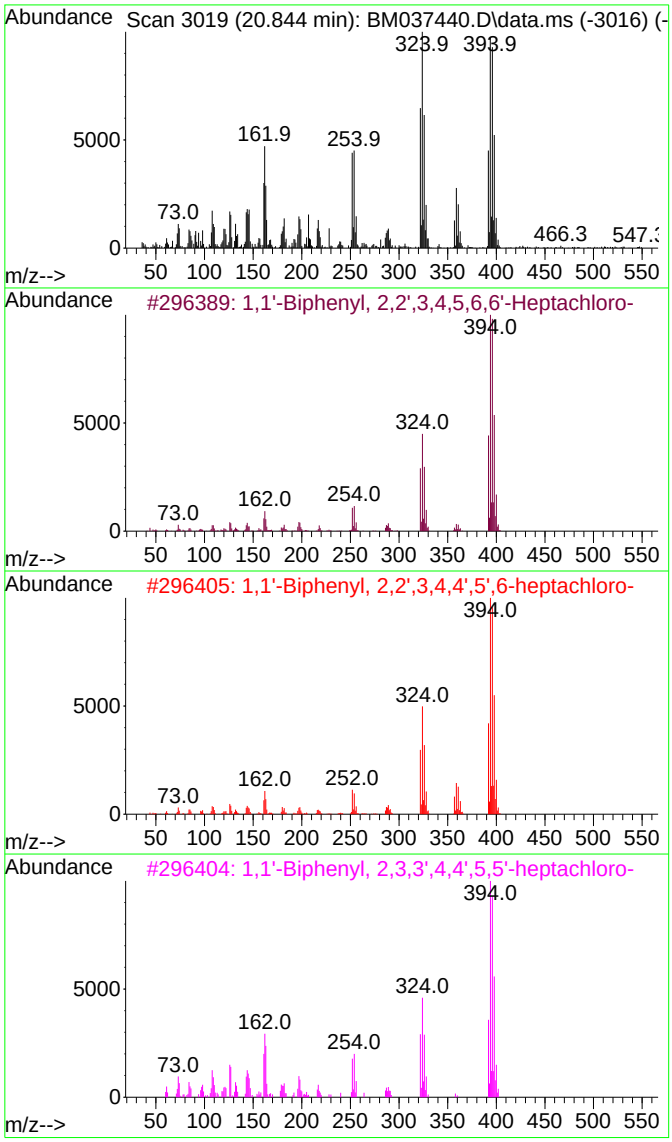
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

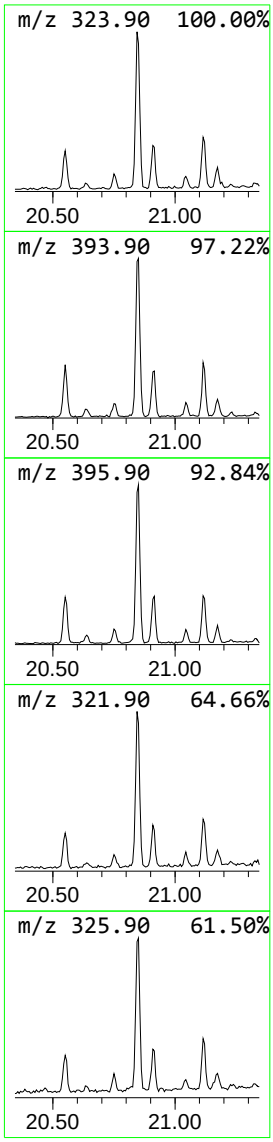
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1,1'-Biphenyl, 2,2',3,4,5,6... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.844	2.33 ng	634985	Chrysene-d12	21.386

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,4,5,6,6'-H...	392	C12H3Cl7	074472-49-4	99
2	1,1'-Biphenyl, 2,2',3,4,4',5',6-...	392	C12H3Cl7	052663-69-1	99
3	1,1'-Biphenyl, 2,3,3',4,4',5,5'-...	392	C12H3Cl7	039635-31-9	99
4	2,2',3,4',5,6,6'-Heptachloro-1,1...	392	C12H3Cl7	074487-85-7	99
5	1,1'-Biphenyl, 2,2',3,3',4,4',5-...	392	C12H3Cl7	035065-30-6	99





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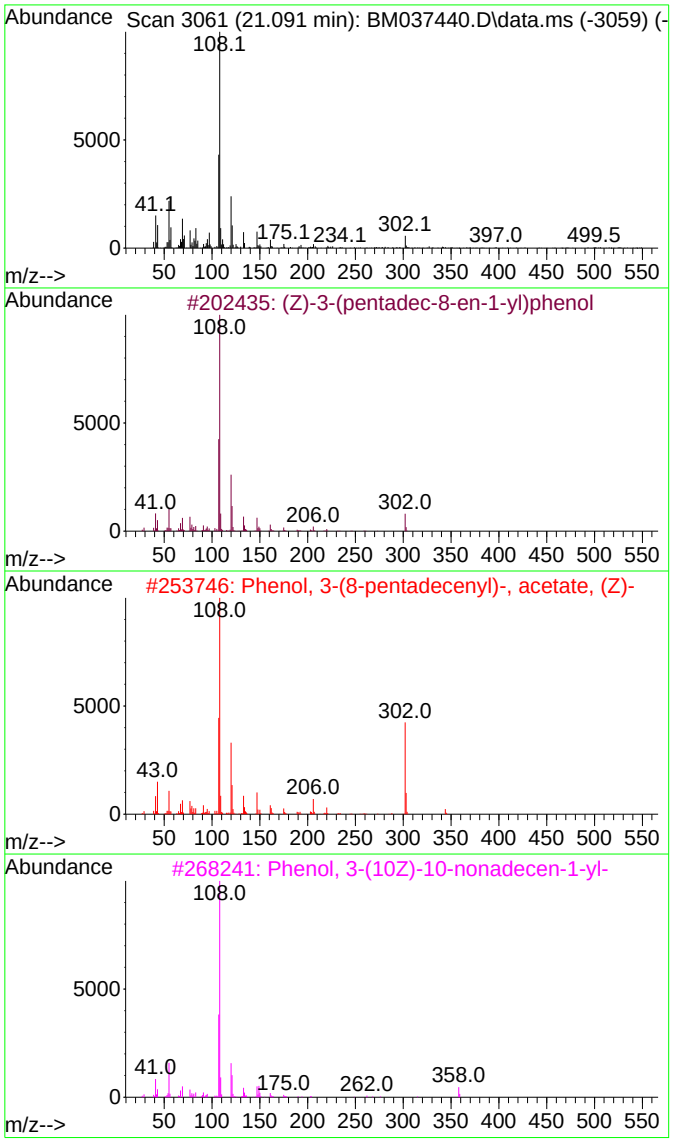
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

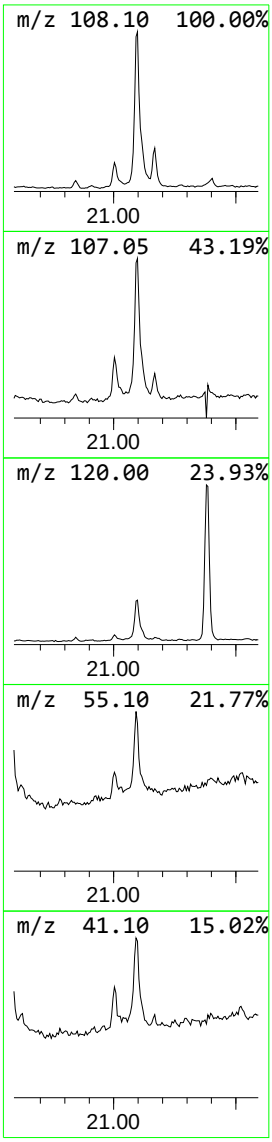
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (Z)-3-(pentadec-8-en-1-yl)p... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.092	3.82 ng	1043540	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	94
2	Phenol, 3-(8-pentadecenyl)-, ace...	344	C23H36O2	111047-34-8	94
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	74
5	Phenol, 3-docosyl-	402	C28H50O	498557-21-4	64





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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-04

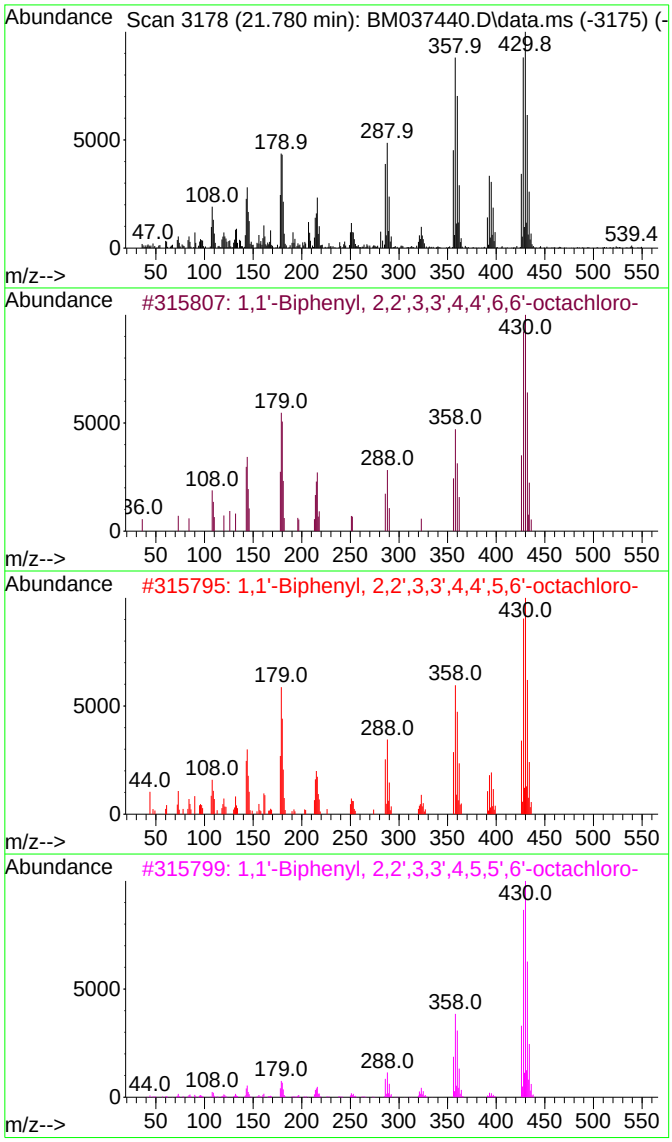
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

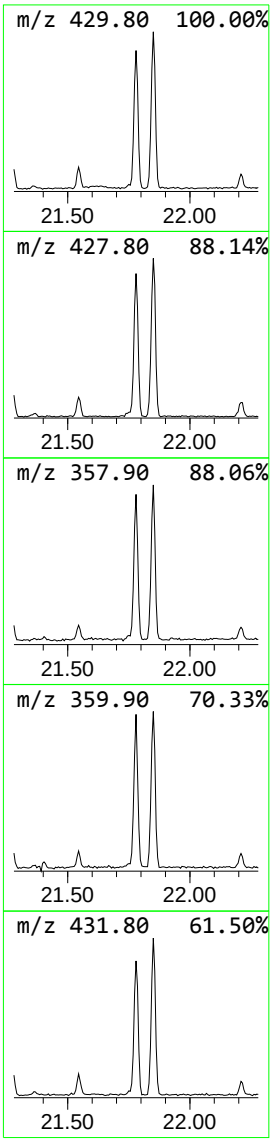
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.780	3.42 ng	932310	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	042740-50-1	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	052663-78-2	99
5	1,1'-Biphenyl, 2,2',3,3',4,5,6,6...	426	C12H2Cl8	052663-73-7	99





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ALS Vial : 9 Sample Multiplier: 1

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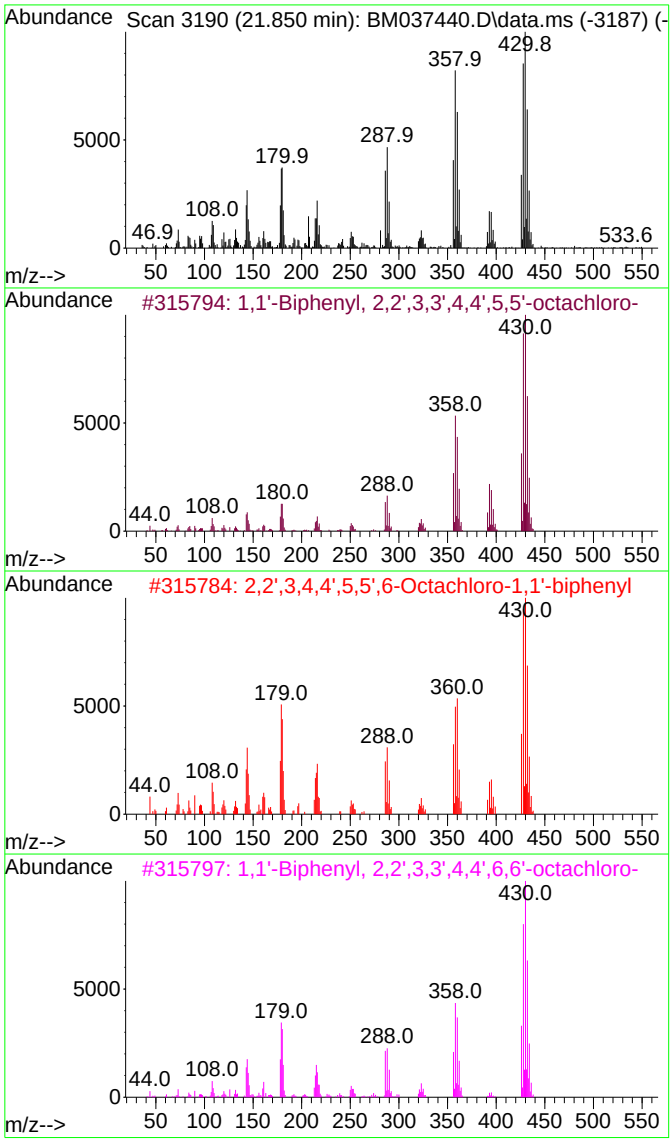
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

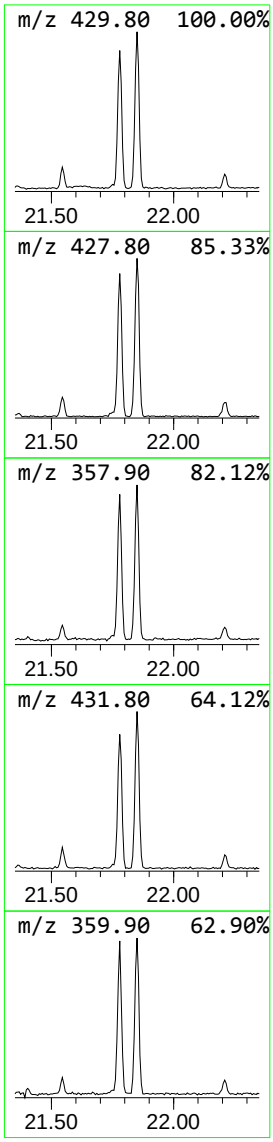
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.850	2.79 ng	760380	Chrysene-d12	21.386

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	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',6,...	426	C12H2Cl8	033091-17-7	99
4	1,1'-Biphenyl, 2,2',3,3',4,5,6,6...	426	C12H2Cl8	052663-73-7	99
5	1,1'-Biphenyl, 2,2',3,3',5,5',6,...	426	C12H2Cl8	002136-99-4	99





Tentatively Identified Compound (LSC) summary

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone, 4-...	4.934	2.0	ng	159425	1	7.822	1577430	20.0
9H-Pyrido[3,4-b...	14.857	2.1	ng	308986	3	14.463	3004060	20.0
o-Terphenyl	17.851	9.6	ng	1699560	4	17.203	3550730	20.0
n-Hexadecanoic ...	18.098	8.4	ng	1483360	4	17.203	3550730	20.0
Phenol, 2-[(4-h...	18.603	4.3	ng	762864	4	17.203	3550730	20.0
Phenol, 4,4'-me...	19.015	4.2	ng	739594	4	17.203	3550730	20.0
Octadecanoic acid	19.380	2.5	ng	686098	5	21.386	5459270	20.0
1,1'-Biphenyl, ...	20.844	2.3	ng	634985	5	21.386	5459270	20.0
(Z)-3-(pentadec...	21.092	3.8	ng	1043540	5	21.386	5459270	20.0
1,1'-Biphenyl, ...	21.780	3.4	ng	932310	5	21.386	5459270	20.0
1,1'-Biphenyl, ...	21.850	2.8	ng	760380	5	21.386	5459270	20.0