

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121521\
 Data File : BP008413.D
 Acq On : 15 Dec 2021 12:59
 Operator : CG/JU
 Sample : M5076-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-19-G

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_P\Methods\8270E-BP121421.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP008413.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.875	316	321	329	rVB	188721	262547	15.07%	1.883%
2	5.346	395	401	421	rBV	530580	878619	50.43%	6.300%
3	5.946	499	503	512	rVB	19074	28141	1.62%	0.202%
4	6.940	666	672	698	rBV	457857	916478	52.61%	6.572%
5	7.740	802	808	819	rBV	199971	326574	18.75%	2.342%
6	8.911	998	1007	1025	rBV	305877	596272	34.23%	4.276%
7	10.540	1276	1284	1296	rBV	207104	403771	23.18%	2.895%
8	13.016	1698	1705	1718	rBV	723930	1137174	65.27%	8.154%
9	13.722	1820	1825	1830	rBV4	9944	19818	1.14%	0.142%
10	14.398	1933	1940	1947	rBV	319622	497158	28.54%	3.565%
11	14.463	1947	1951	1959	rVB2	14665	26277	1.51%	0.188%
12	15.904	2187	2196	2215	rBV	454507	844758	48.49%	6.058%
13	16.828	2348	2353	2359	rBV	27173	42162	2.42%	0.302%
14	16.975	2374	2378	2390	rVB	48501	78837	4.53%	0.565%
15	17.157	2403	2409	2413	rBV	379816	580515	33.32%	4.163%
16	17.204	2413	2417	2424	rVV	211710	318594	18.29%	2.285%
17	17.298	2428	2433	2440	rVB	29069	51601	2.96%	0.370%
18	17.745	2504	2509	2518	rVB	59881	89686	5.15%	0.643%
19	17.881	2527	2532	2540	rVB3	33351	70994	4.07%	0.509%
20	18.034	2553	2558	2563	rBV2	61587	100993	5.80%	0.724%
21	18.081	2563	2566	2573	rVV	77293	113363	6.51%	0.813%
22	18.216	2584	2589	2593	rVV2	60782	106273	6.10%	0.762%
23	18.257	2593	2596	2602	rVB	55880	76717	4.40%	0.550%
24	18.522	2635	2641	2645	rBV	41199	69794	4.01%	0.500%
25	18.575	2647	2650	2658	rVB2	52990	83356	4.78%	0.598%
26	18.928	2706	2710	2714	rBV	30726	44190	2.54%	0.317%
27	19.075	2729	2735	2739	rBV2	37115	59918	3.44%	0.430%
28	19.181	2748	2753	2760	rBV	187277	251033	14.41%	1.800%
29	19.533	2806	2813	2820	rBV	283684	400796	23.01%	2.874%
30	19.728	2841	2846	2858	rBV	1363438	1742188	100.00%	12.493%
31	19.869	2865	2870	2876	rVB4	20175	41528	2.38%	0.298%
32	20.004	2887	2893	2900	rBV4	29232	79245	4.55%	0.568%
33	20.181	2920	2923	2929	rBV	37866	42990	2.47%	0.308%
34	20.316	2942	2946	2949	rBV	32022	41915	2.41%	0.301%
35	20.363	2950	2954	2957	rVV2	34188	43601	2.50%	0.313%
36	20.763	3019	3022	3026	rVB	38193	43312	2.49%	0.311%

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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_P\Methods\8270E-BP121421.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	20.975	3056	3058	3066	rVB3	73230	109303	6.27%	0.784%
38	21.263	3101	3107	3111	rBV2	622571	945784	54.29%	6.782%
39	21.304	3111	3114	3122	rVV	124946	172582	9.91%	1.238%
40	21.380	3122	3127	3143	rVV	877830	1381860	79.32%	9.909%
41	23.633	3504	3510	3518	rVB2	426769	824697	47.34%	5.914%

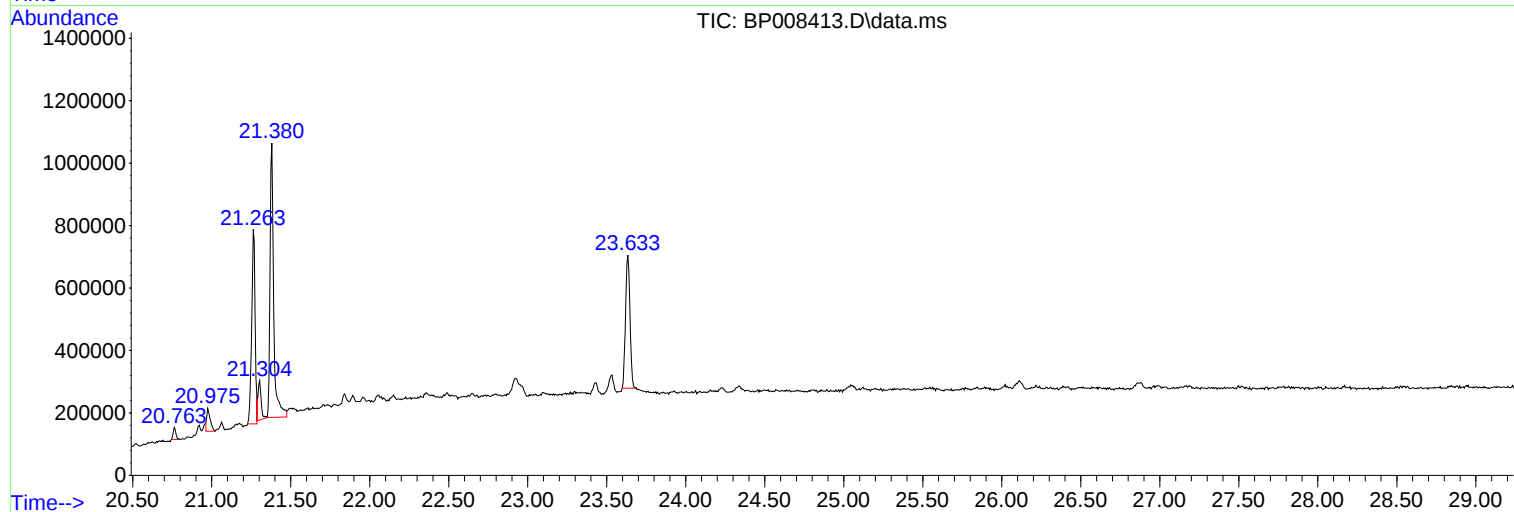
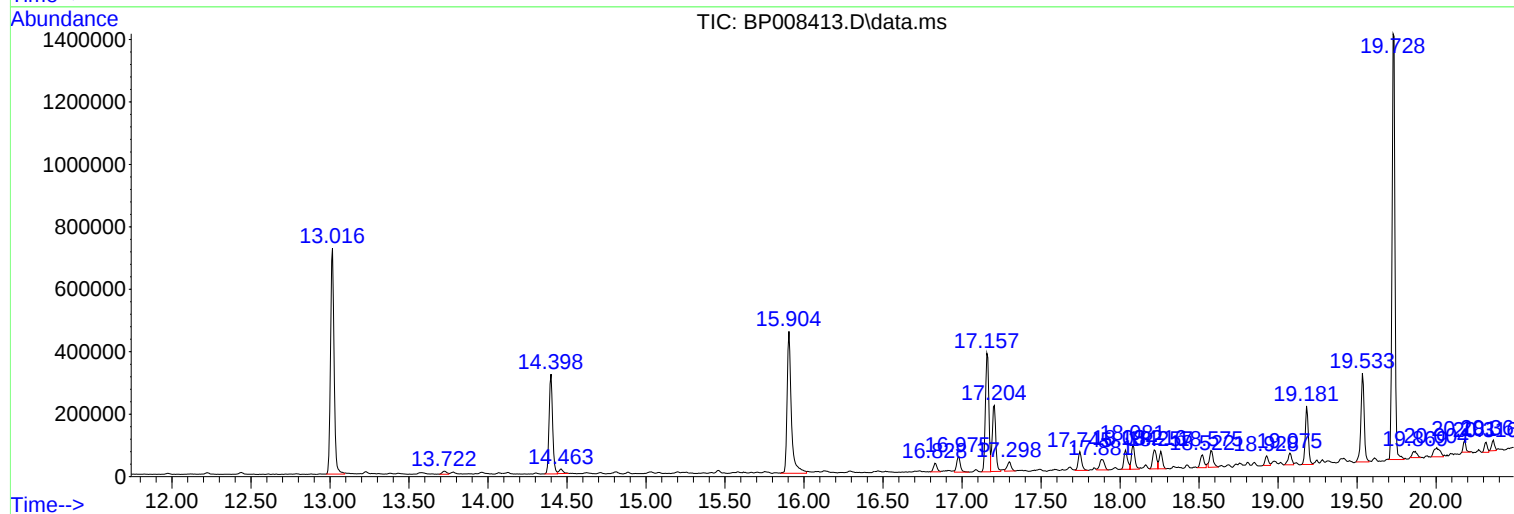
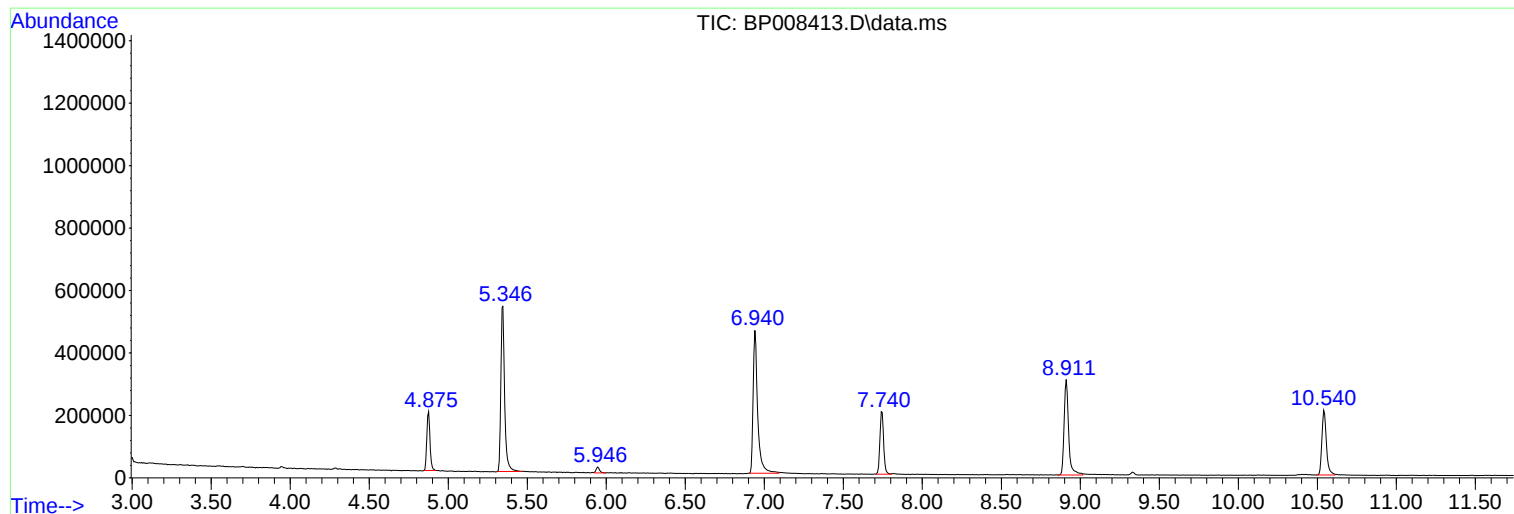
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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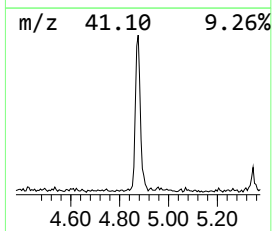
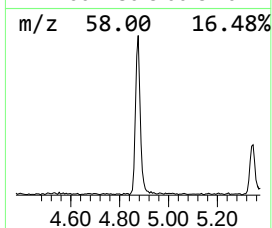
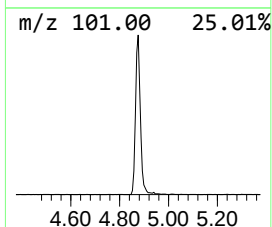
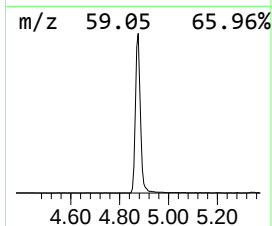
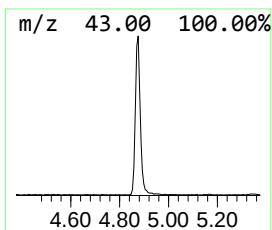
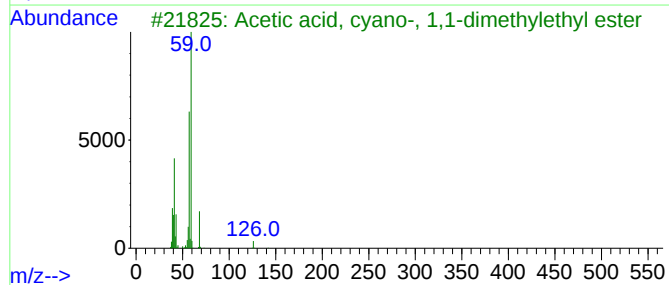
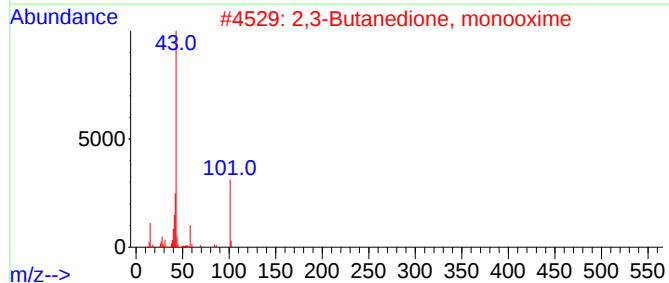
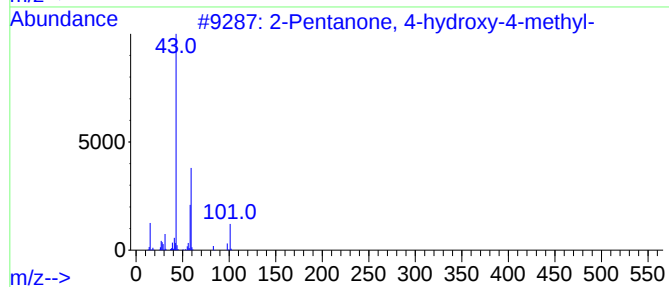
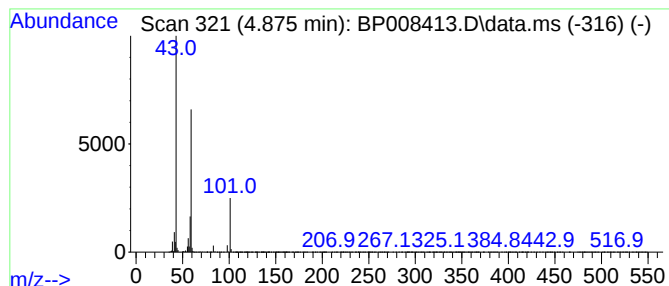
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.875	16.08 ng	262547	1,4-Dichlorobenzene-d4	7.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72		
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17		
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		
5	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9		



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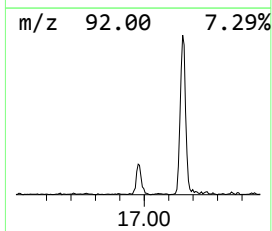
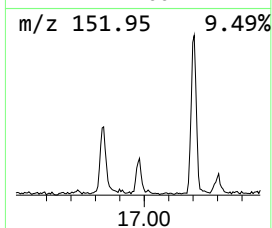
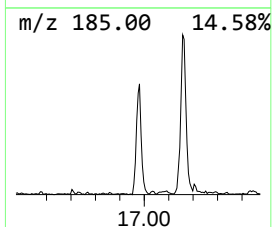
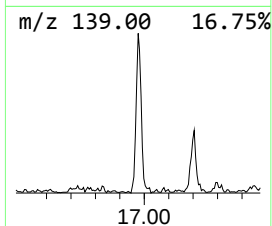
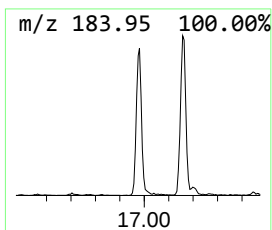
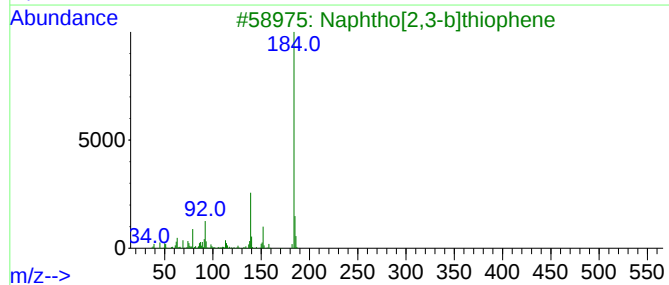
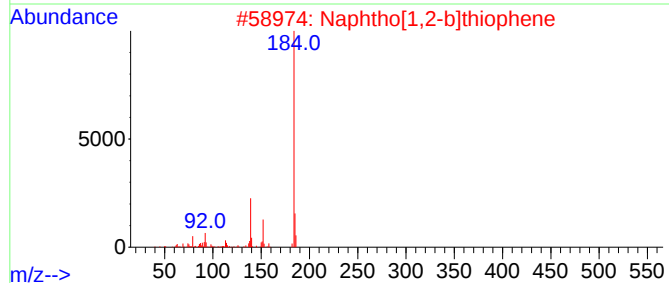
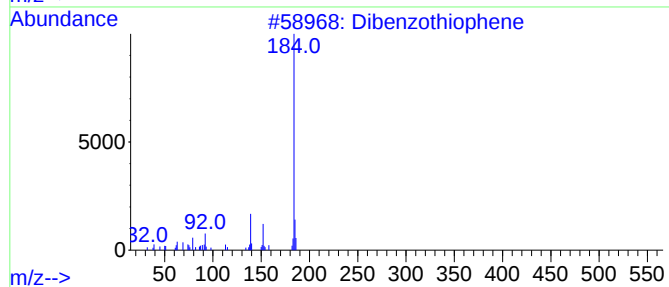
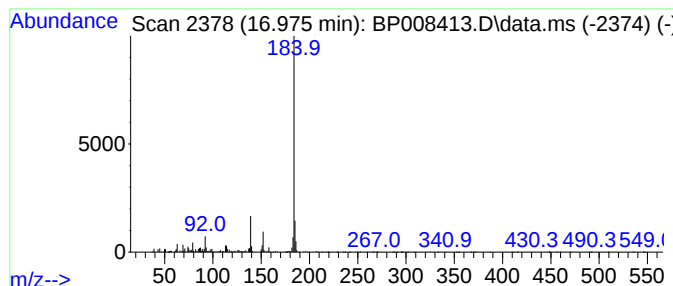
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Dibenzothiophene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.975	2.72 ng	78837	Phenanthrene-d10	17.157

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	97
2			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	93
3			Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	93
4			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	91
5			Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	90



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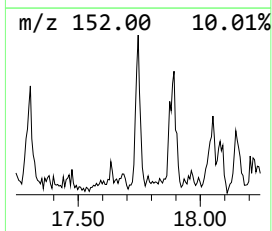
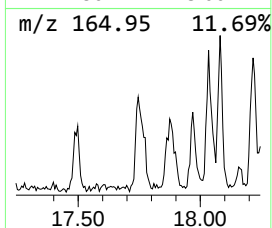
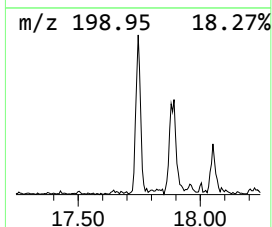
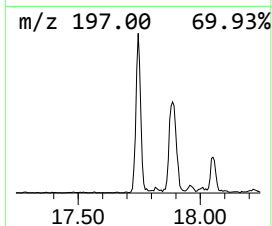
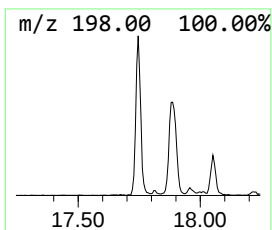
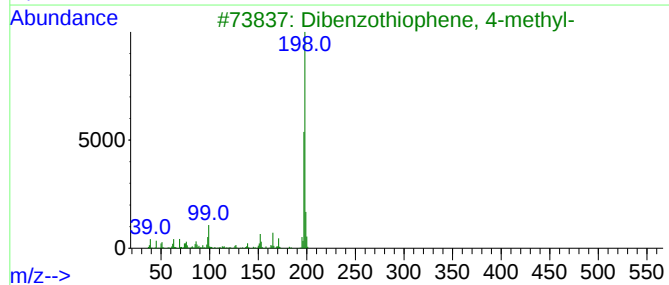
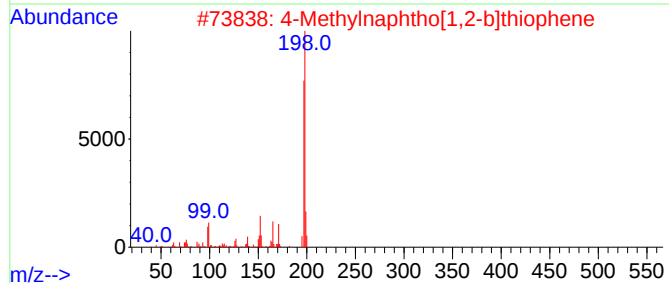
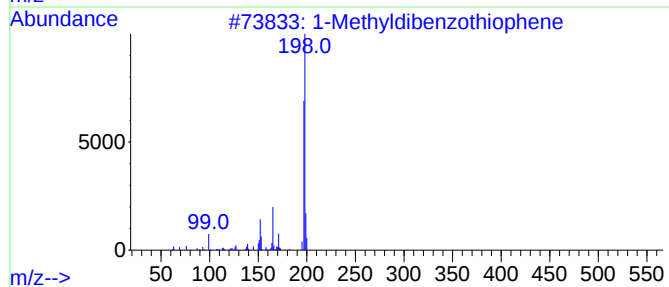
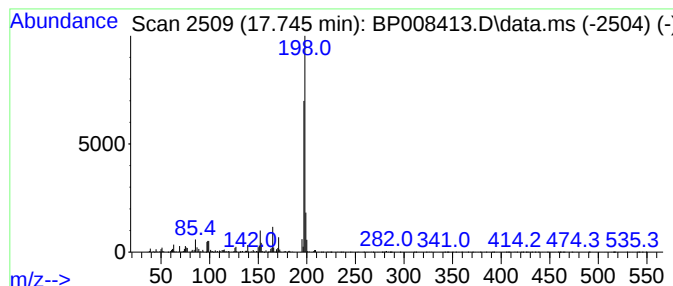
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Methyldibenzothiophene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.745	3.09 ng	89686	Phenanthrene-d10	17.157

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methyldibenzothiophene	198	C13H10S	031317-07-4	97
2			4-Methylnaphtho[1,2-b]thiophene	198	C13H10S	067388-11-8	96
3			Dibenzothiophene, 4-methyl-	198	C13H10S	007372-88-5	95
4			Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	95
5			2-Methyldibenzothiophene	198	C13H10S	1000383-55-1	93



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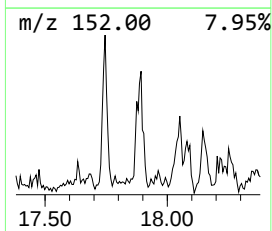
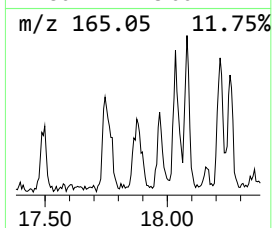
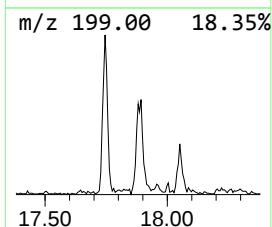
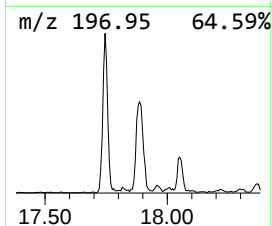
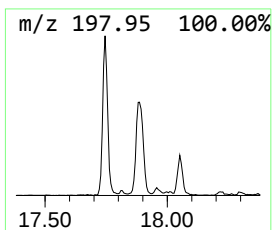
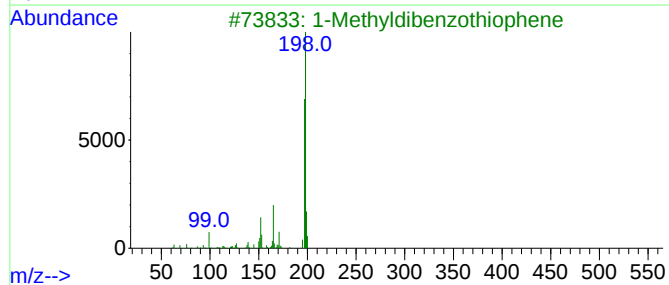
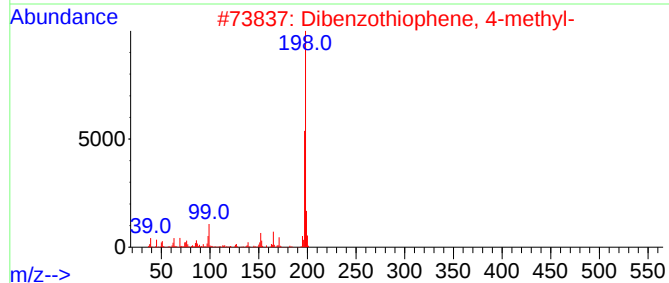
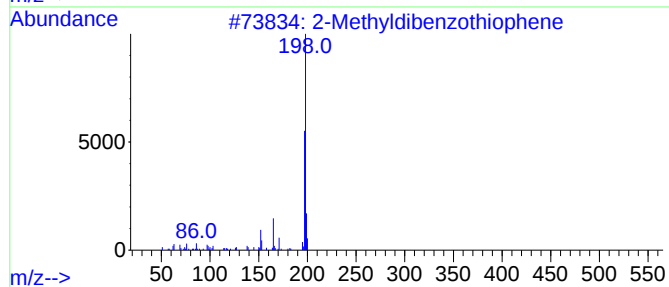
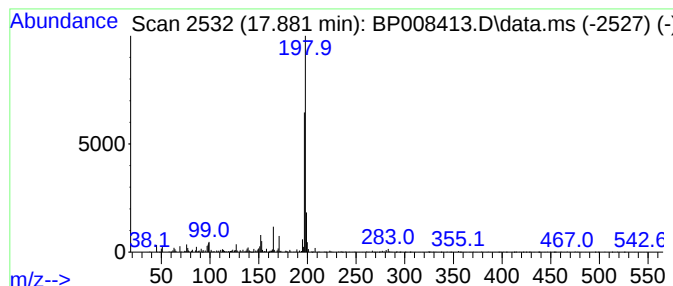
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 2-Methyldibenzothiophene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.881	2.45 ng	70994	Phenanthrene-d10	17.157

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methyldibenzothiophene	198	C13H10S	1000383-55-1	96
2			Dibenzothiophene, 4-methyl-	198	C13H10S	007372-88-5	93
3			1-Methyldibenzothiophene	198	C13H10S	031317-07-4	91
4			Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	90
5			4-Methylnaphtho[1,2-b]thiophene	198	C13H10S	067388-11-8	90



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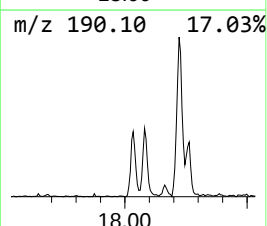
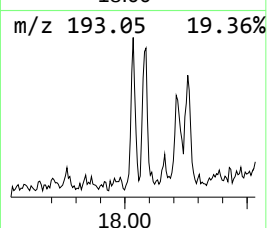
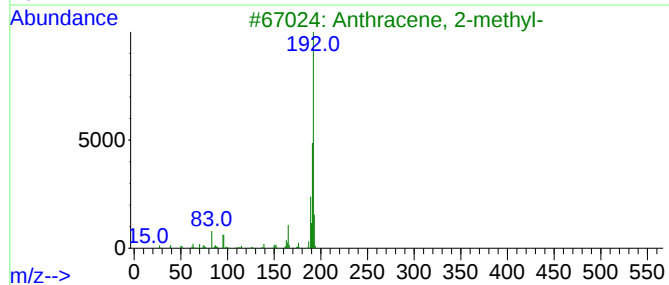
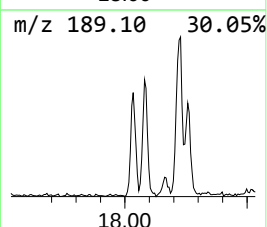
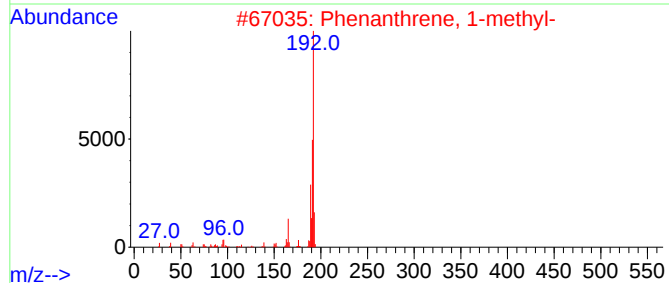
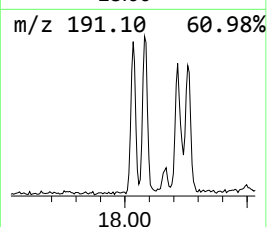
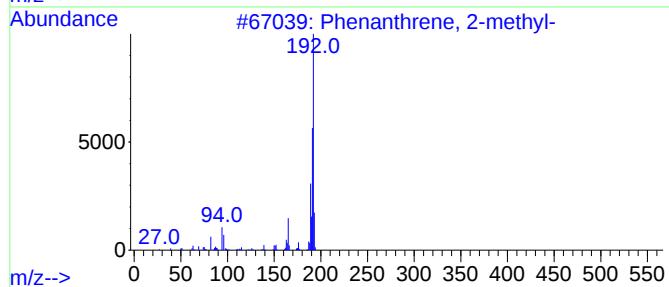
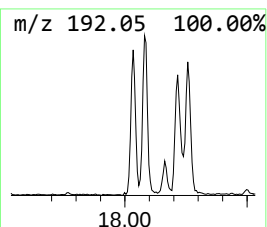
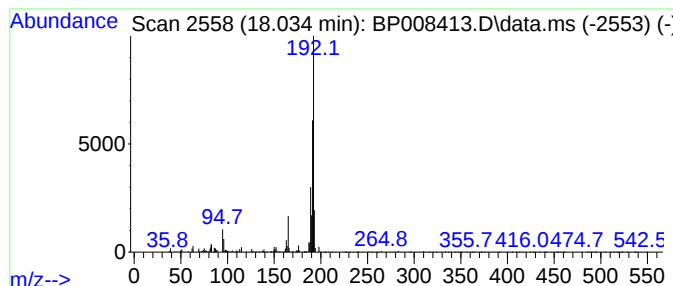
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.034	3.48 ng	100993	Phenanthrene-d10	17.157

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121521\
Data File : BP008413.D
Acq On : 15 Dec 2021 12:59
Operator : CG/JU
Sample : M5076-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-19-G

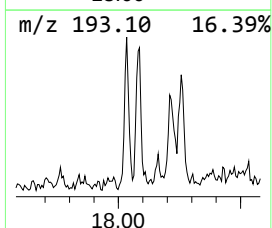
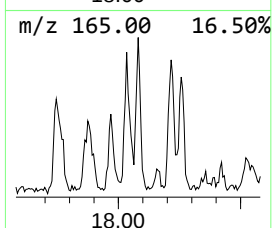
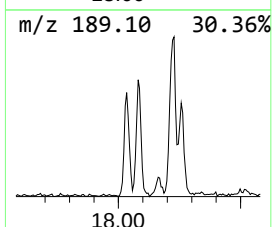
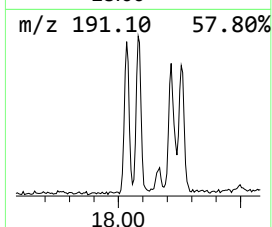
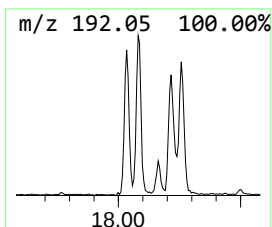
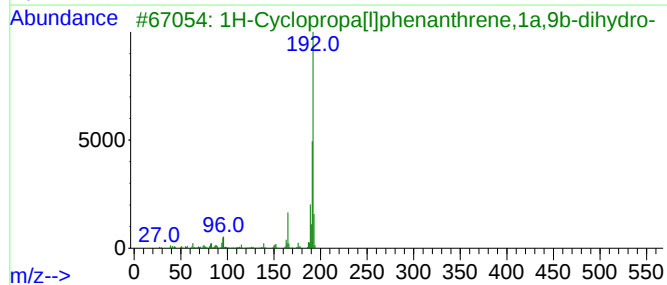
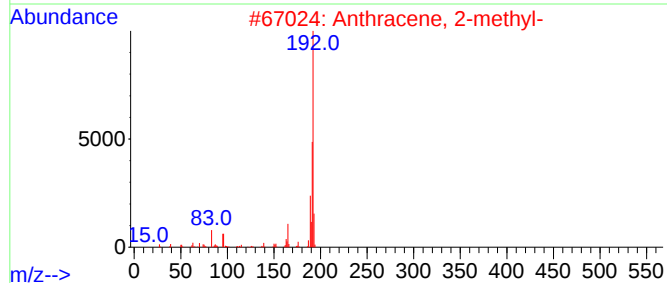
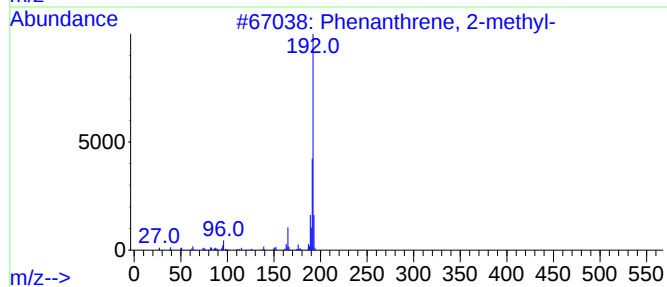
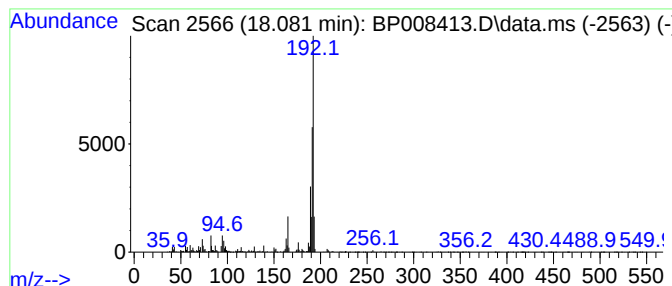
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.081	3.91 ng	113363	Phenanthrene-d10	17.157

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
3		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	93
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93
5		Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121521\
Data File : BP008413.D
Acq On : 15 Dec 2021 12:59
Operator : CG/JU
Sample : M5076-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

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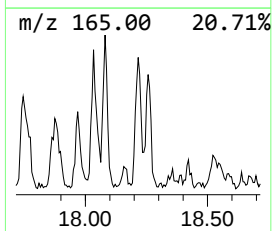
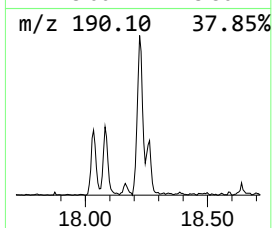
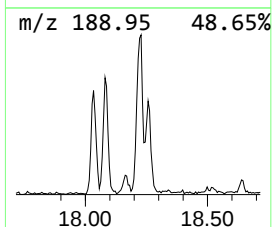
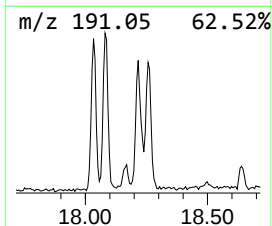
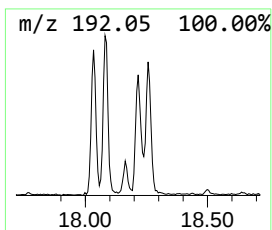
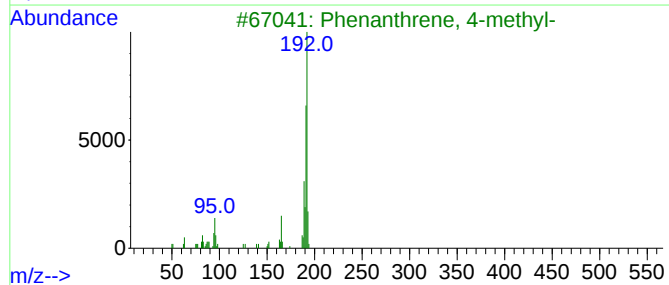
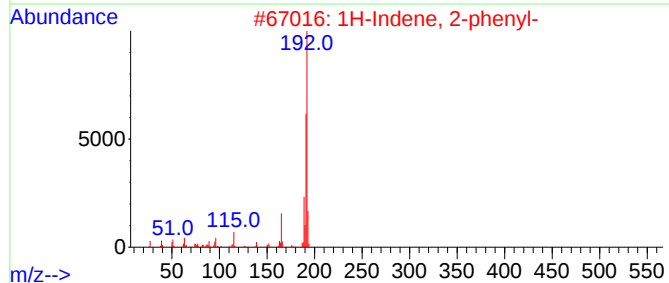
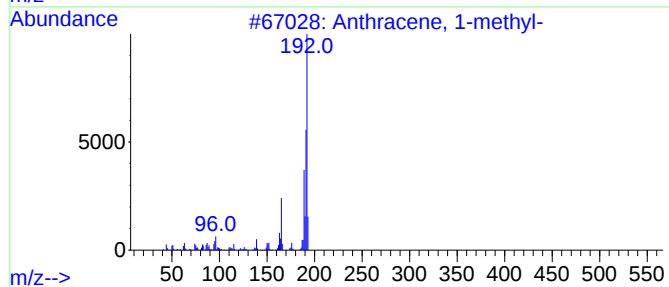
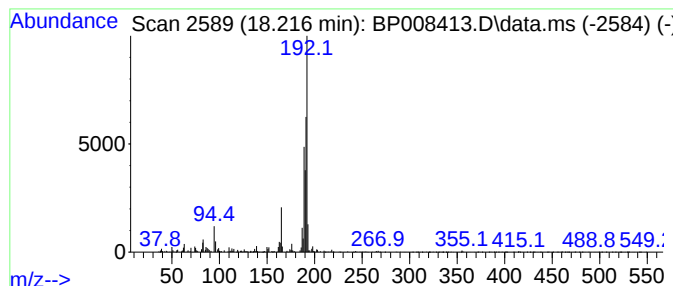
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Anthracene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	3.66 ng	106273	Phenanthrene-d10	17.157

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	64
2			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	64
3			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	60
4			1a,9b-Dihydro-1H-cyclopropa[a]an...	192	C15H12	006109-24-6	58
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121521\
Data File : BP008413.D
Acq On : 15 Dec 2021 12:59
Operator : CG/JU
Sample : M5076-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

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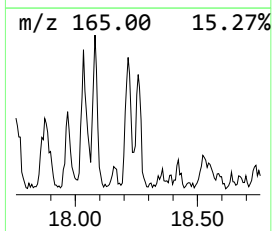
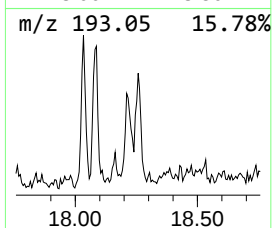
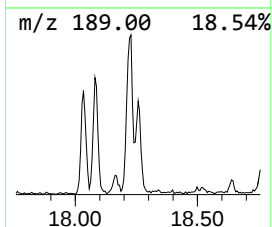
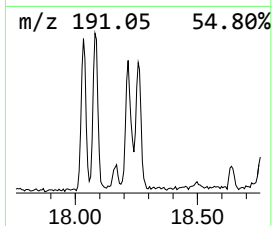
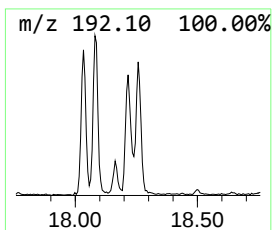
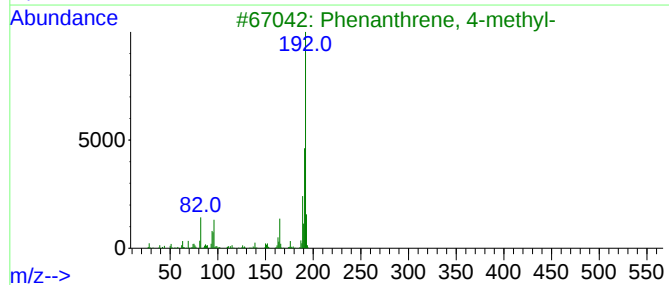
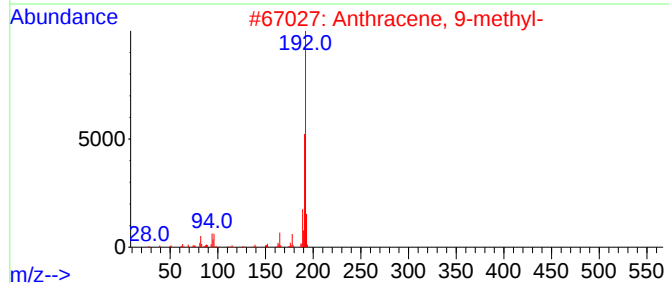
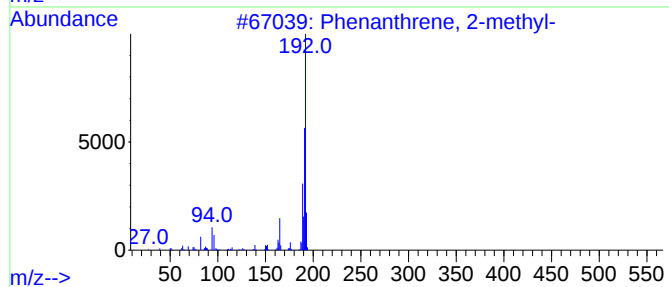
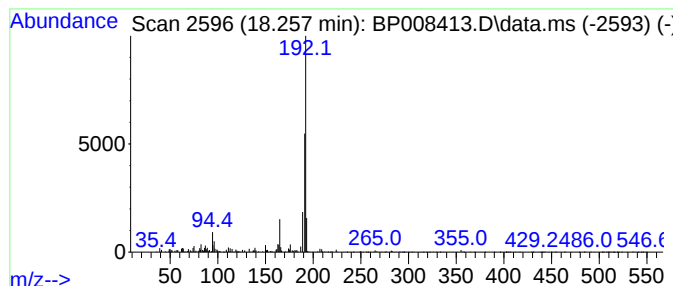
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Anthracene, 9-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.257	2.64 ng	76717	Phenanthrene-d10	17.157

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
2		Anthracene, 9-methyl-	192	C15H12	000779-02-2	87
3		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	87
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	87
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121521\
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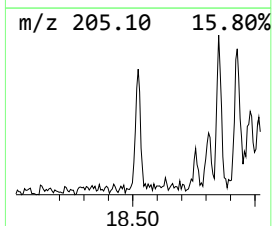
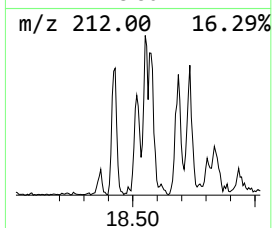
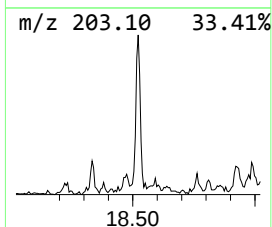
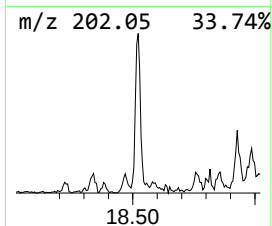
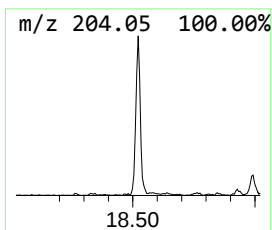
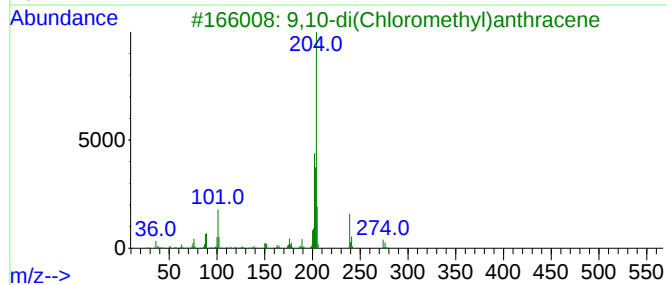
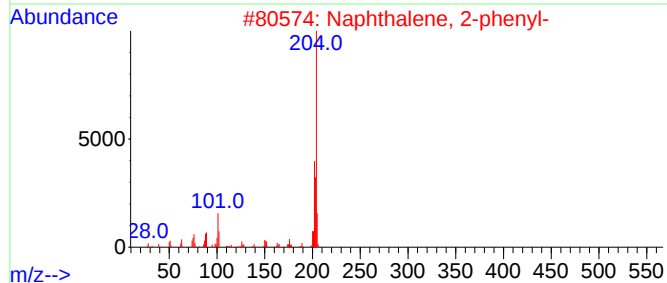
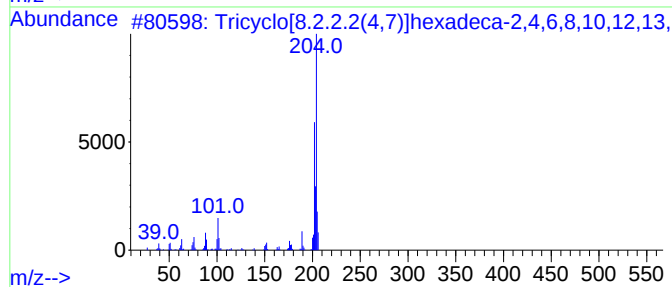
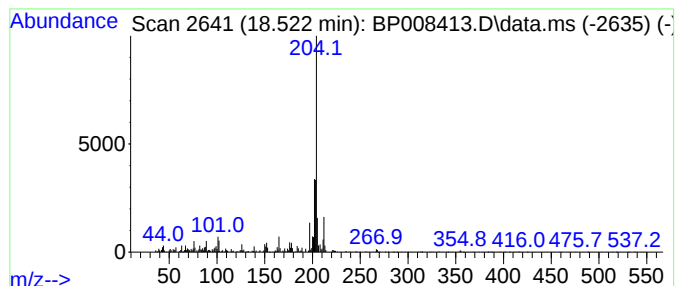
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tricyclo[8.2.2.2(4,7)]hexad... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.522	2.40 ng	69794	Phenanthrene-d10	17.157	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	64
2	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	64
3	9,10-di(Chloromethyl)anthracene	274	C16H12Cl2	010387-13-0	64
4	3,4-Biphenyldicarbonitrile	204	C14H8N2	004128-63-6	52
5	[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121521\
Data File : BP008413.D
Acq On : 15 Dec 2021 12:59
Operator : CG/JU
Sample : M5076-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

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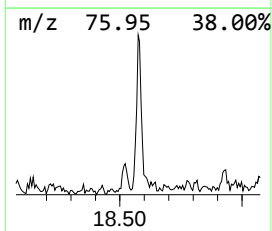
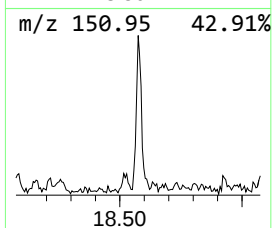
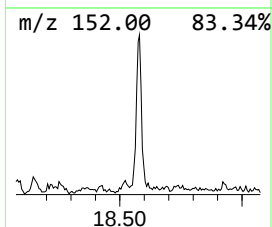
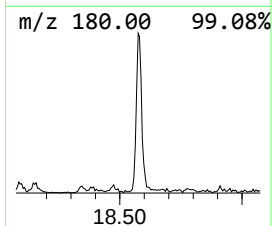
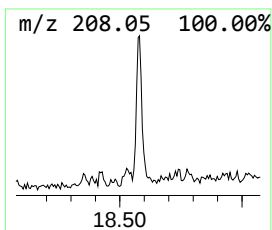
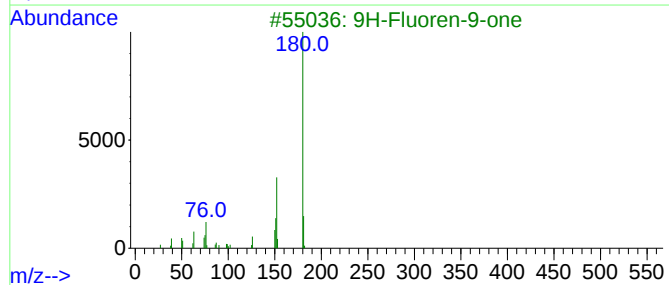
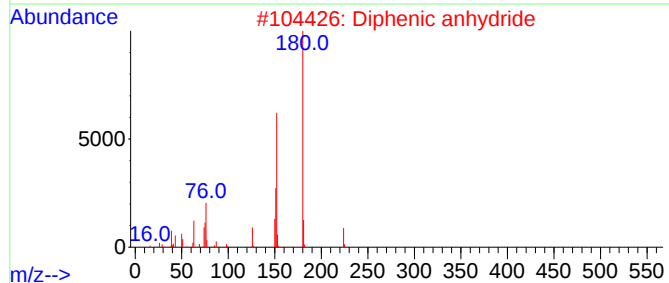
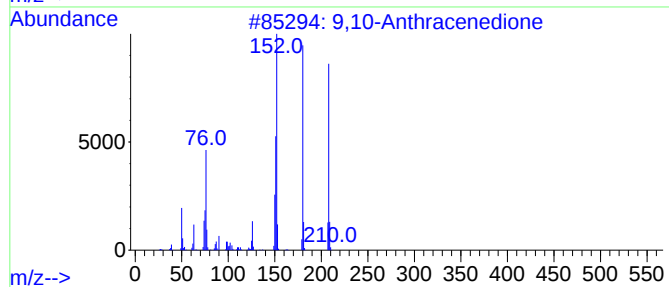
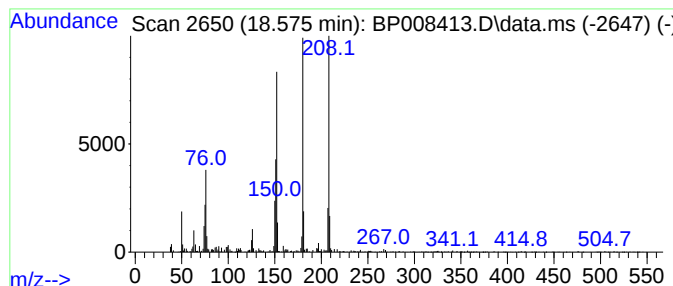
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9,10-Anthracenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.575	2.87 ng	83356	Phenanthrene-d10	17.157

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,10-Anthracenedione	208	C14H8O2	000084-65-1	96
2		Diphenic anhydride	224	C14H8O3	006050-13-1	64
3		9H-Fluoren-9-one	180	C13H8O	000486-25-9	64
4		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	64
5		Benzo[h]cinnoline	180	C12H8N2	000230-31-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121521\
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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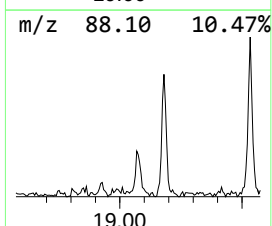
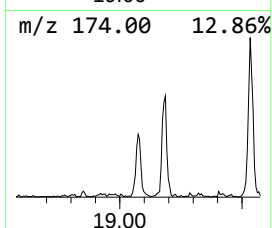
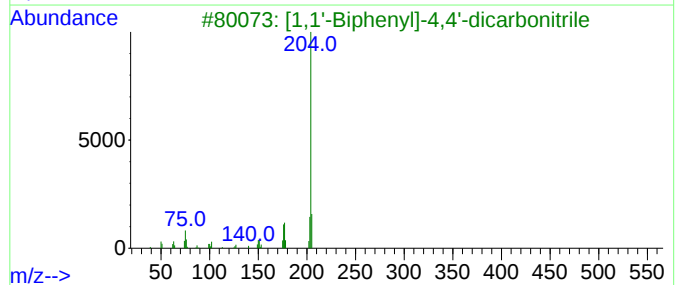
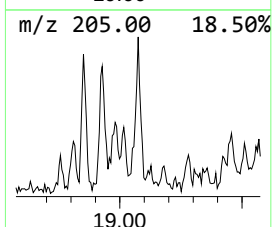
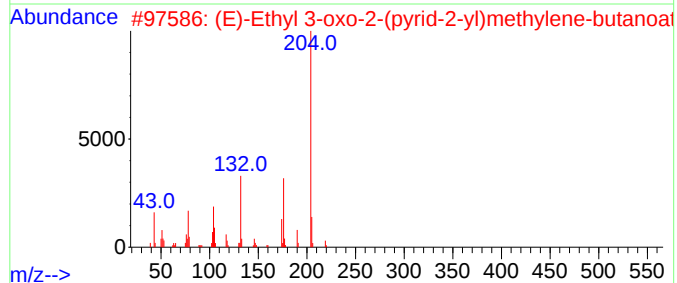
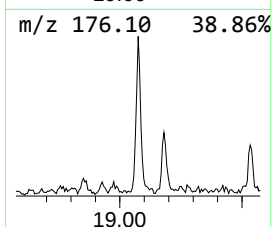
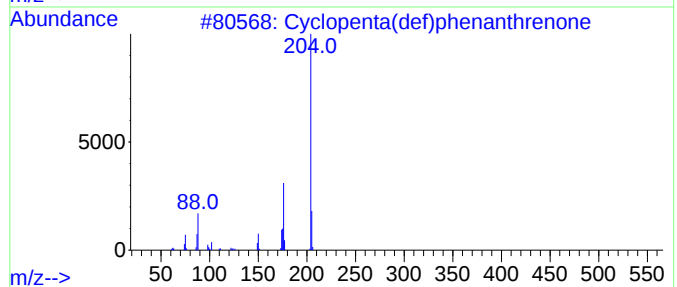
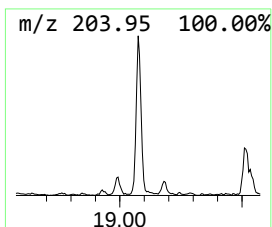
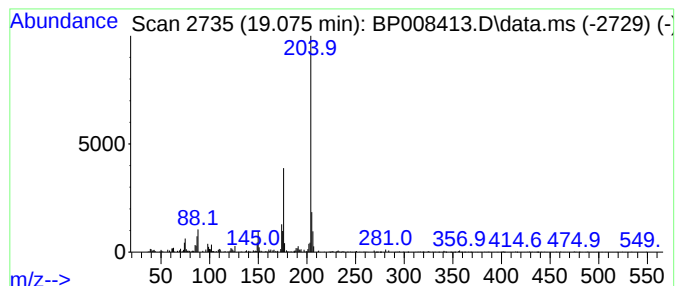
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Cyclopenta(def)phenanthrenone Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.075	2.06 ng	59918	Phenanthrene-d10	17.157

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	91
2			(E)-Ethyl 3-oxo-2-(pyrid-2-yl)me...	219	C12H13NO3	1000147-59-7	59
3			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	50
4			6-Phenyl-2-thiouracil	204	C10H8N2OS	005590-94-3	50
5			9-Acridinecarbonitrile	204	C14H8N2	005326-19-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121521\
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Sample : M5076-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

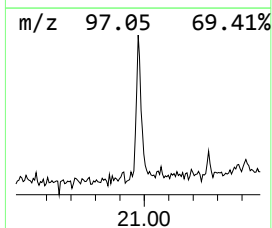
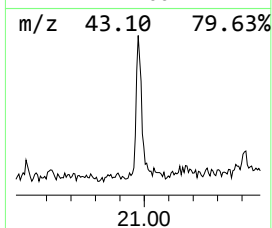
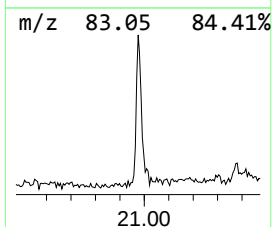
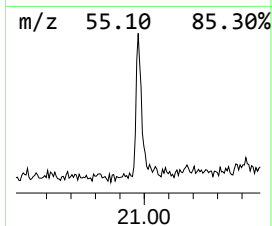
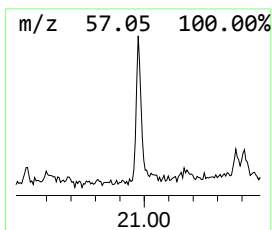
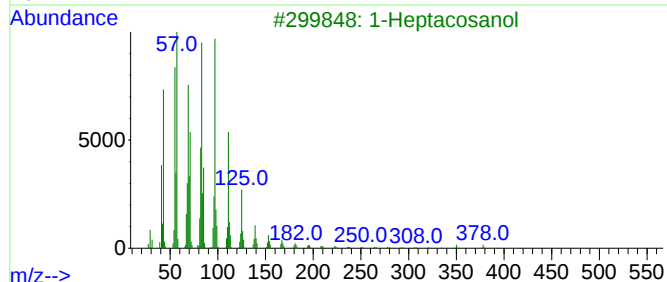
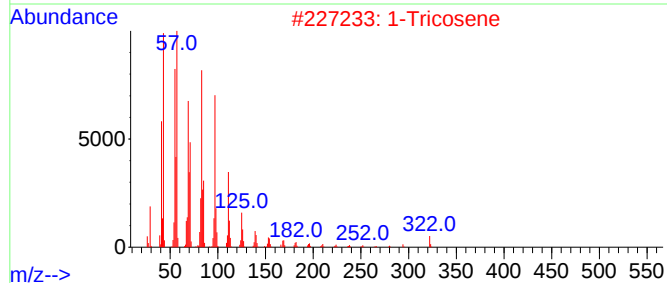
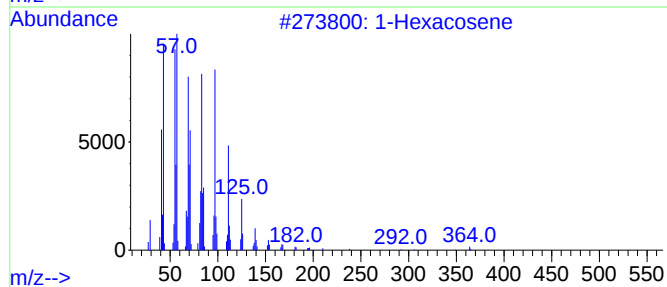
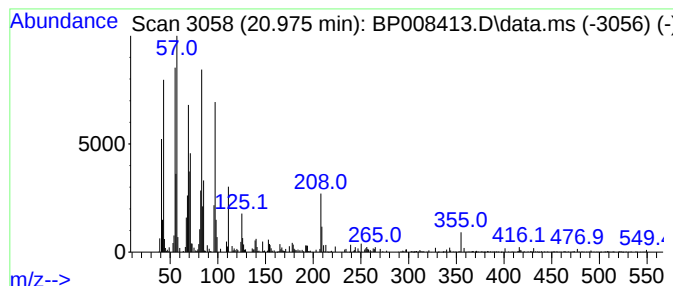
Instrument :
BNA_P
ClientSampleId :
SB-19-G

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.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-Hexacosene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.975	2.31 ng	109303	Chrysene-d12	21.263	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene	364	C26H52	018835-33-1	87
2	1-Tricosene	322	C23H46	018835-32-0	87
3	1-Heptacosanol	396	C27H56O	002004-39-9	87
4	Cyclotetracosane	336	C24H48	000297-03-0	83
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121521\
Data File : BP008413.D
Acq On : 15 Dec 2021 12:59
Operator : CG/JU
Sample : M5076-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-19-G

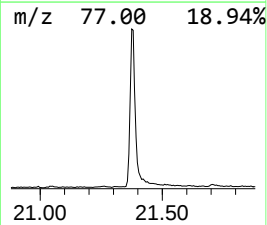
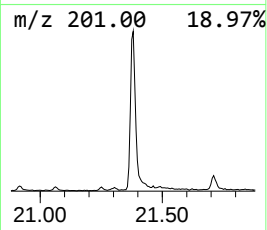
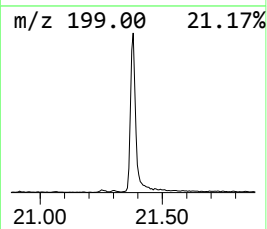
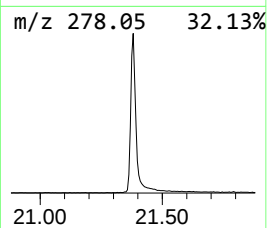
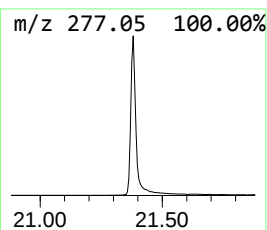
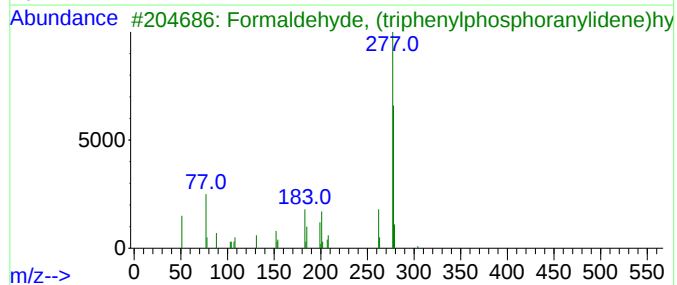
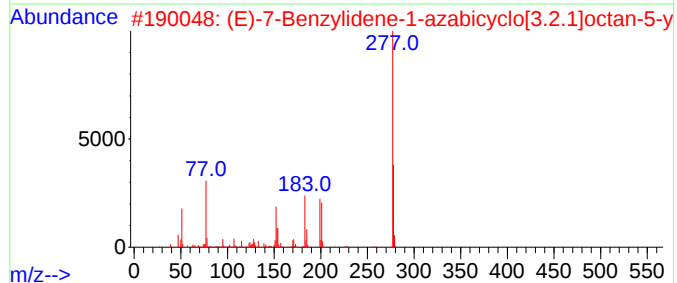
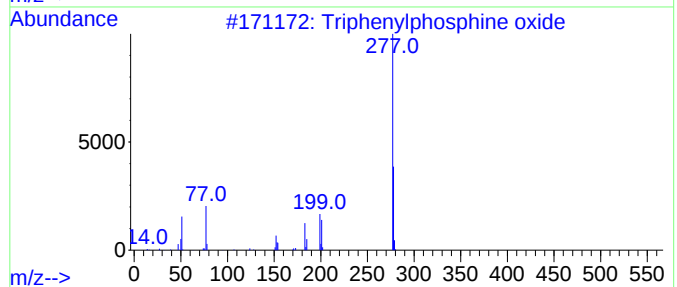
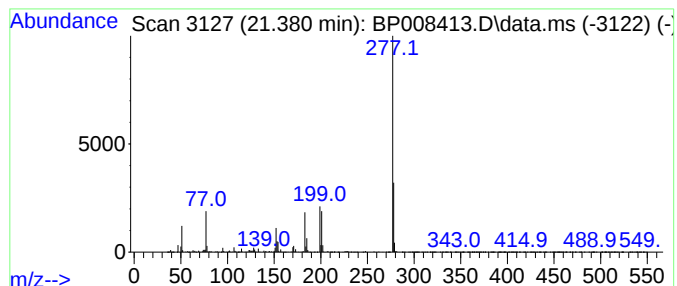
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Triphenylphosphine oxide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.381	29.22 ng	1381860	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	96
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19NO3S	1000483-83-4	62
3			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	42
4			3-(3,4-Dichlorophenyl)-2-thioxo-...	277	C9H5Cl2NOS2	017046-34-3	38
5			5-Methyl-8-phenylfuro[3',2':5,6]...	277	C16H11N3O2	1000460-07-9	28



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121521\
Data File : BP008413.D
Acq On : 15 Dec 2021 12:59
Operator : CG/JU
Sample : M5076-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-19-G

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.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-...	4.875	16.1	ng	262547	1	7.740	326574	20.0
Dibenzothiophene	16.975	2.7	ng	78837	4	17.157	580515	20.0
1-Methyldibenzo...	17.745	3.1	ng	89686	4	17.157	580515	20.0
2-Methyldibenzo...	17.881	2.5	ng	70994	4	17.157	580515	20.0
Phenanthrene, 2...	18.034	3.5	ng	100993	4	17.157	580515	20.0
Anthracene, 2-m...	18.081	3.9	ng	113363	4	17.157	580515	20.0
Anthracene, 1-m...	18.216	3.7	ng	106273	4	17.157	580515	20.0
Anthracene, 9-m...	18.257	2.6	ng	76717	4	17.157	580515	20.0
Tricyclo[8.2.2....	18.522	2.4	ng	69794	4	17.157	580515	20.0
9,10-Anthracene...	18.575	2.9	ng	83356	4	17.157	580515	20.0
Cyclopenta(def)...	19.075	2.1	ng	59918	4	17.157	580515	20.0
1-Hexacosene	20.975	2.3	ng	109303	5	21.263	945784	20.0
Triphenylphosph...	21.381	29.2	ng	1381860	5	21.263	945784	20.0