

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100925\
 Data File : BP025967.D
 Acq On : 09 Oct 2025 16:27
 Operator : RC/JU
 Sample : Q3313-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETGI-360

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-BP091225.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP025967.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.008	6	12	27	rVB	1173475	1750585	25.88%	2.859%
2	5.378	408	415	441	rBV	2077320	3941577	58.26%	6.437%
3	6.943	673	681	709	rBV	1878662	4125150	60.97%	6.737%
4	7.731	808	815	832	rBV	539676	970296	14.34%	1.585%
5	8.878	1003	1010	1034	rBV	1190211	2549315	37.68%	4.164%
6	10.496	1277	1285	1309	rBV	596743	1386616	20.50%	2.265%
7	12.948	1693	1702	1719	rBV	2680507	5258675	77.73%	8.588%
8	14.342	1930	1939	1946	rBV	1028254	1798576	26.59%	2.937%
9	14.407	1947	1950	1960	rVB2	52937	100497	1.49%	0.164%
10	15.419	2116	2122	2132	rBV3	44611	88804	1.31%	0.145%
11	15.713	2166	2172	2189	rBV	392266	637428	9.42%	1.041%
12	15.854	2189	2196	2215	rBV	2392647	4637962	68.55%	7.575%
13	16.954	2379	2383	2394	rVB	47701	76604	1.13%	0.125%
14	17.142	2408	2415	2419	rBV	1332135	2087126	30.85%	3.409%
15	17.189	2419	2423	2429	rVB	861099	1215221	17.96%	1.985%
16	17.289	2434	2440	2448	rVB	115564	199580	2.95%	0.326%
17	17.578	2482	2489	2500	rBV2	77792	206094	3.05%	0.337%
18	18.066	2564	2572	2589	rBV2	869834	1748361	25.84%	2.855%
19	18.183	2589	2592	2597	rVV	50734	76270	1.13%	0.125%
20	18.248	2597	2603	2607	rVV2	185623	341605	5.05%	0.558%
21	18.289	2607	2610	2616	rVB2	77474	124426	1.84%	0.203%
22	18.460	2634	2639	2645	rVB7	61969	105885	1.57%	0.173%
23	18.566	2652	2657	2662	rBV	73541	109303	1.62%	0.179%
24	18.625	2663	2667	2675	rVB2	90149	161619	2.39%	0.264%
25	19.001	2726	2731	2735	rBV	97134	164131	2.43%	0.268%
26	19.160	2752	2758	2762	rBV2	78183	144890	2.14%	0.237%
27	19.272	2770	2777	2784	rBV	2271177	2989939	44.19%	4.883%
28	19.401	2794	2799	2810	rBV3	195407	418353	6.18%	0.683%
29	19.613	2831	2835	2836	rBV2	283205	316739	4.68%	0.517%
30	19.648	2836	2841	2851	rVB	1657476	2542358	37.58%	4.152%
31	19.736	2852	2856	2863	rVB2	81401	130037	1.92%	0.212%
32	19.860	2871	2877	2884	rBV	4949239	6765349	100.00%	11.049%
33	19.995	2893	2900	2906	rVB3	85247	233093	3.45%	0.381%
34	20.172	2927	2930	2937	rVB2	252235	400848	5.93%	0.655%
35	20.272	2943	2947	2951	rBV	137928	187949	2.78%	0.307%
36	20.330	2953	2957	2961	rBV2	127223	205076	3.03%	0.335%

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Smoothing : ON

Filtering: 5

Sampling : 1

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Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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37	20.483	2980	2983	2988	rBV	106553	135356	2.00%	0.221%
38	20.536	2988	2992	2996	rBV	105759	154159	2.28%	0.252%
39	20.707	3018	3021	3025	rBV3	84241	95342	1.41%	0.156%
40	20.971	3062	3066	3073	rBV	158543	237308	3.51%	0.388%
41	21.154	3093	3097	3100	rBV2	163345	225066	3.33%	0.368%
42	21.189	3100	3103	3107	rVB	119525	145935	2.16%	0.238%
43	21.236	3107	3111	3122	rBV4	610705	1074720	15.89%	1.755%
44	21.324	3123	3126	3131	rVB	125283	177789	2.63%	0.290%
45	21.583	3162	3170	3175	rBV2	1533085	3765146	55.65%	6.149%
46	21.630	3175	3178	3192	rVB	831505	1369156	20.24%	2.236%
47	23.865	3552	3558	3566	rBV	491084	1526055	22.56%	2.492%
48	24.612	3679	3685	3698	rVB	375942	958230	14.16%	1.565%
49	24.760	3705	3710	3722	rVB	471476	1209798	17.88%	1.976%
50	24.912	3729	3736	3745	rBV	715805	1959166	28.96%	3.200%

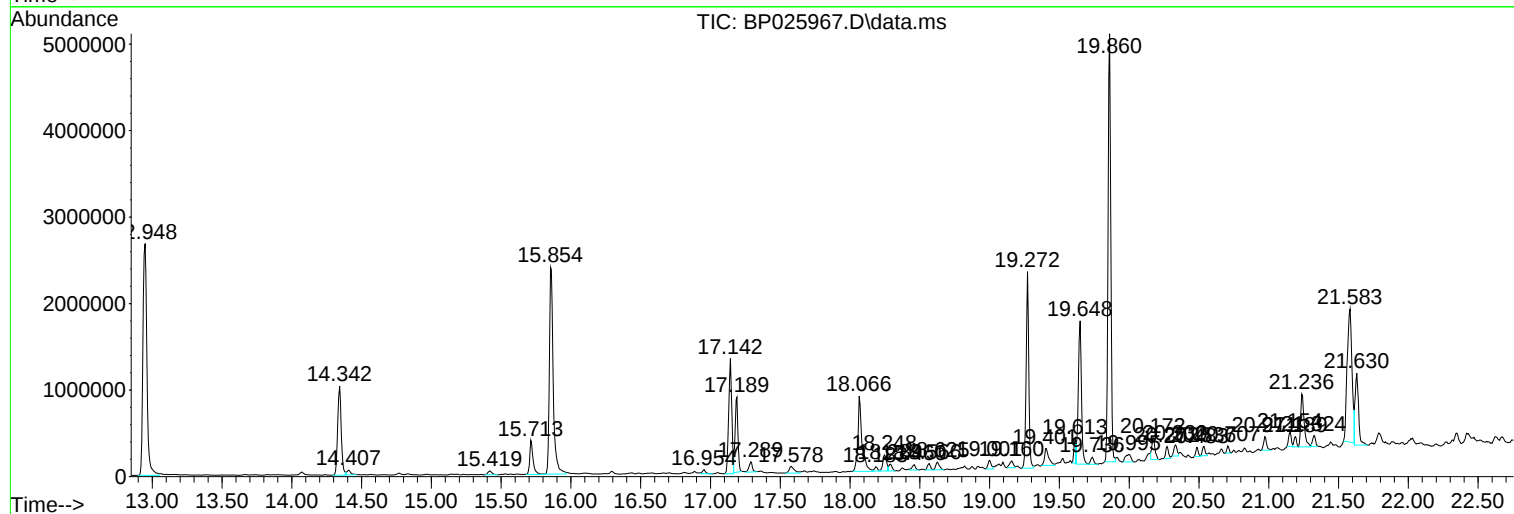
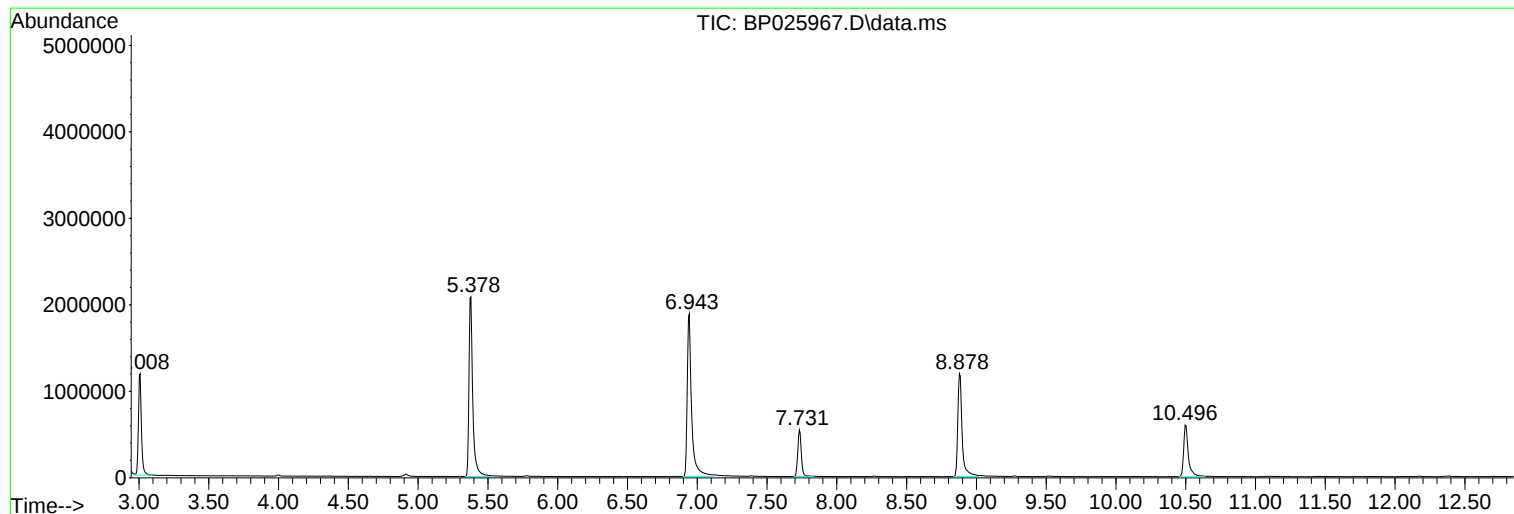
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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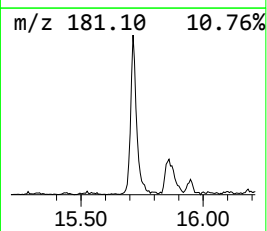
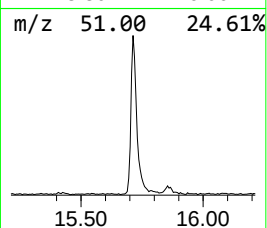
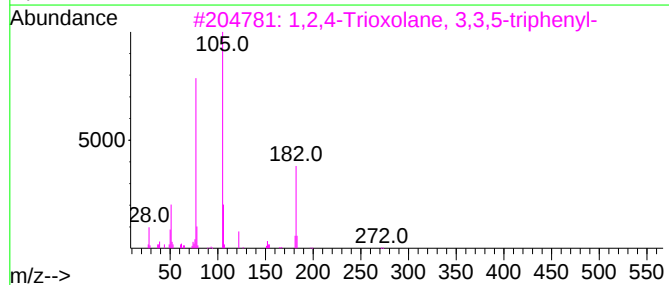
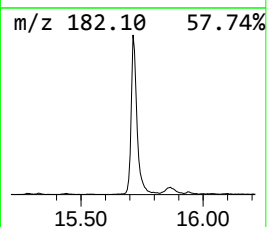
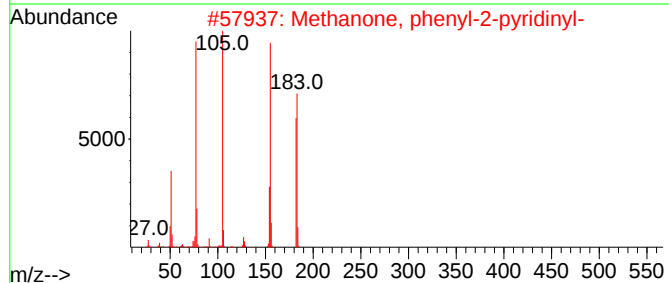
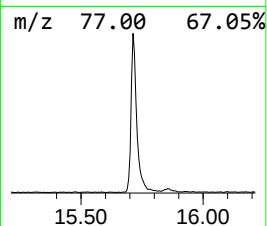
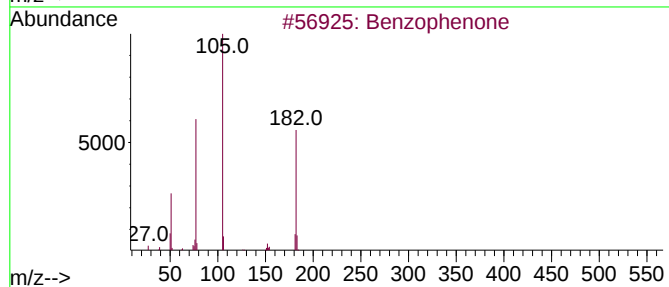
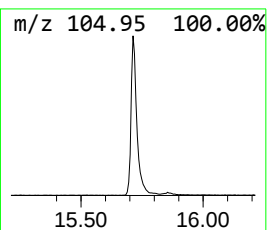
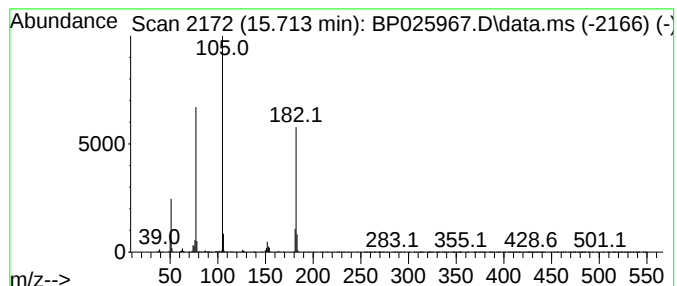
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.713	7.09 ng	637428	Acenaphthene-d10	14.342

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Methanone, phenyl-2-pyridinyl-	183	C12H9NO	000091-02-1	72
3			1,2,4-Trioxolane, 3,3,5-triphenyl-	304	C20H16O3	023246-12-0	64
4			1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	64
5			Azobenzene	182	C12H10N2	000103-33-3	42



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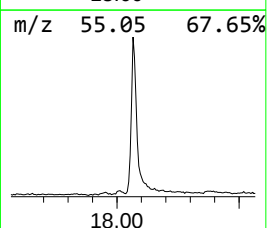
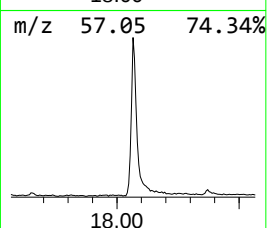
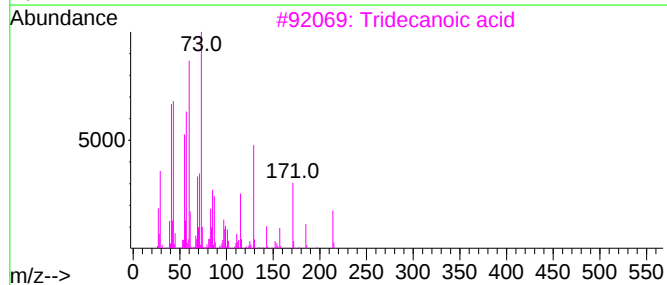
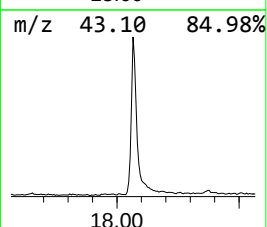
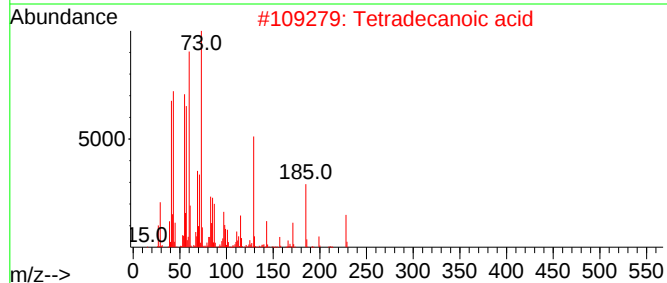
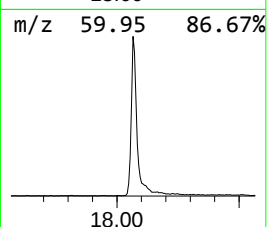
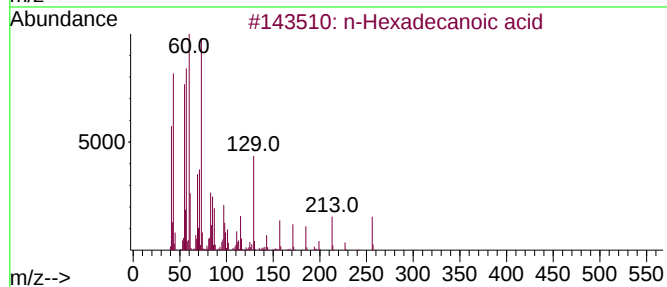
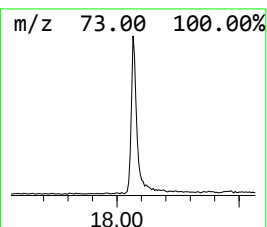
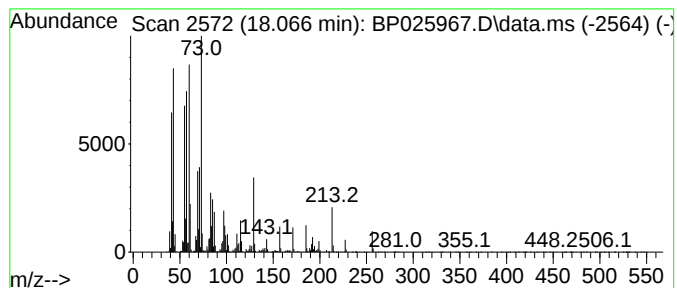
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.066	16.75 ng	1748360	Phenanthrene-d10		17.142	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid		228	C14H28O2	000544-63-8	94
3	Tridecanoic acid		214	C13H26O2	000638-53-9	93
4	Pentadecanoic acid		242	C15H30O2	001002-84-2	87
5	8-Methylnonanoic acid		172	C10H20O2	005963-14-4	55



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Misc :
ALS Vial : 8 Sample Multiplier: 1

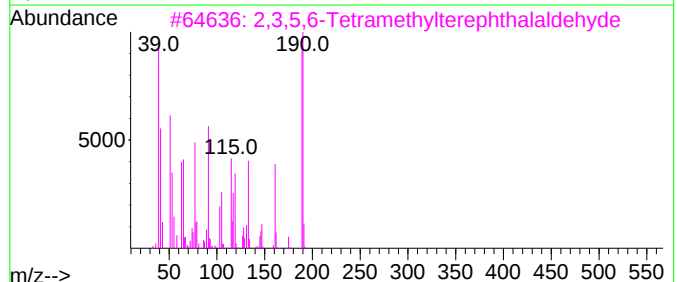
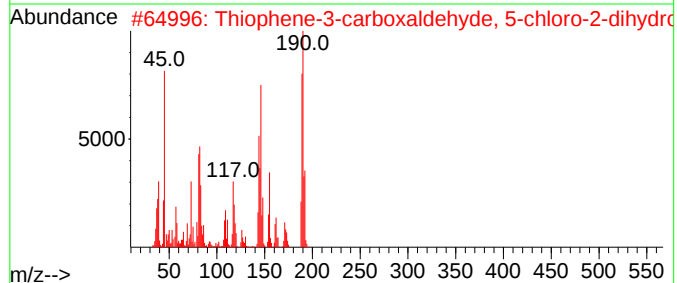
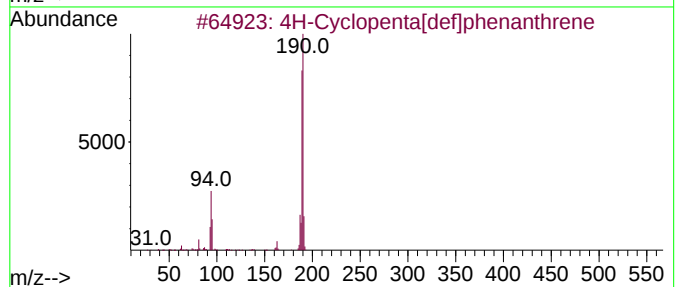
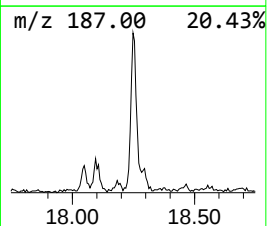
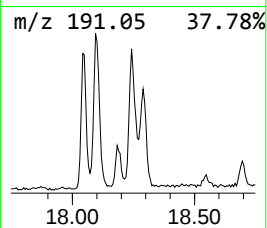
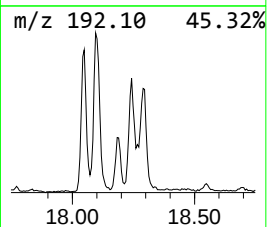
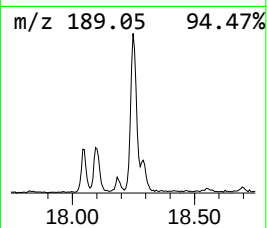
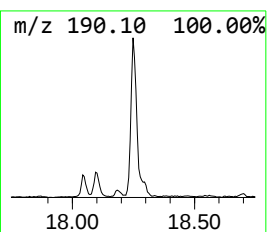
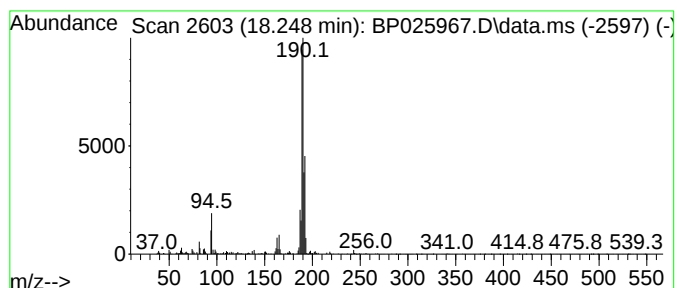
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.248	3.27 ng	341605	Phenanthrene-d10		17.142	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	94
2	Thiophene-3-carboxaldehyde, 5-ch...		190	C5H4BClO3S	036155-87-0	50
3	2,3,5,6-Tetramethylterephthalald...		190	C12H14O2	007072-01-7	47
4	6H-Cyclobuta[jk]phenanthrene		190	C15H10	083469-43-6	45
5	2,2'-Bis(4,5-dimethylimidazole)		190	C10H14N4	069286-06-2	42



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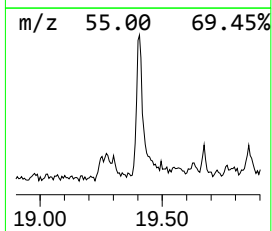
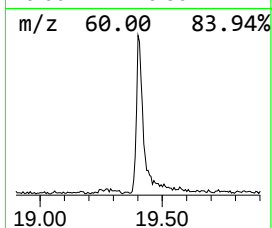
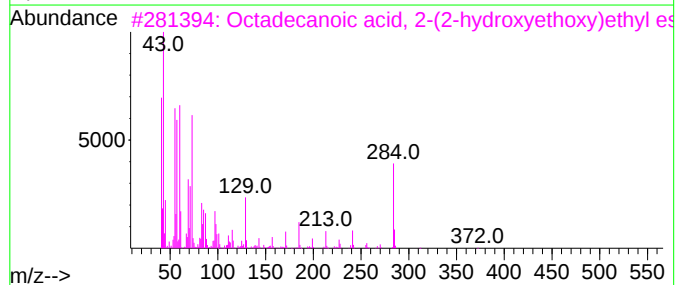
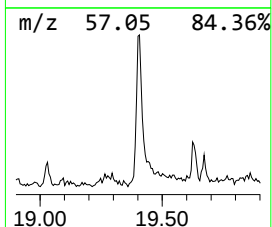
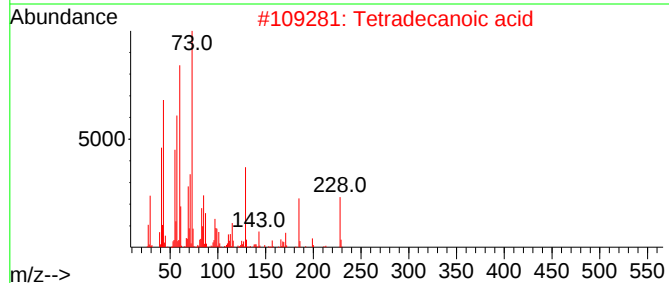
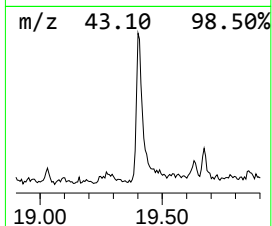
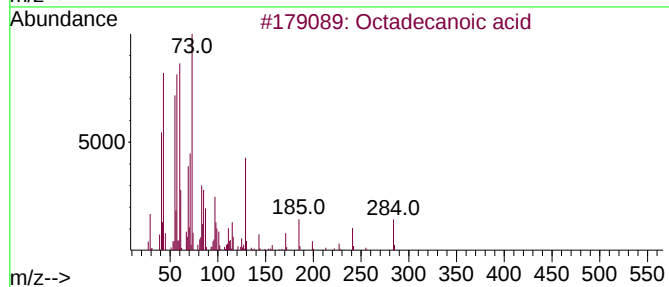
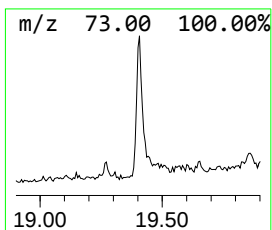
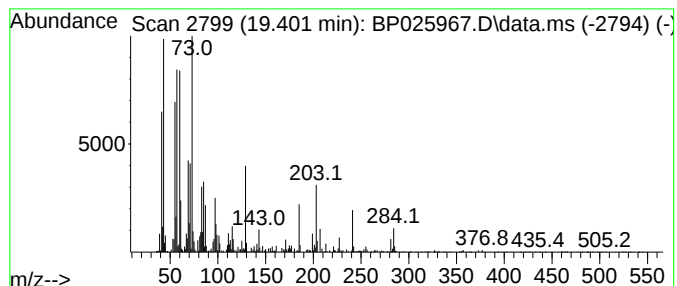
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.401	2.22 ng	418353	Chrysene-d12	21.583	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
3	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	64
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	64
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	62



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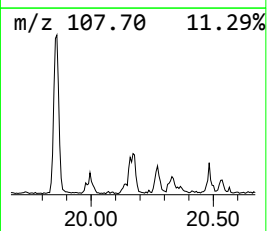
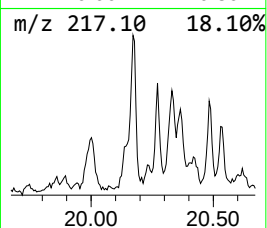
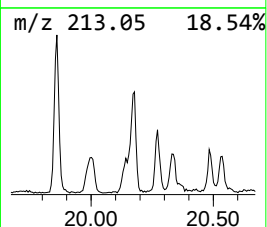
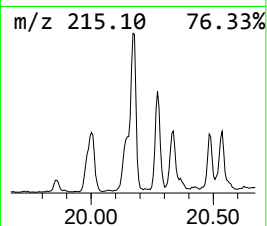
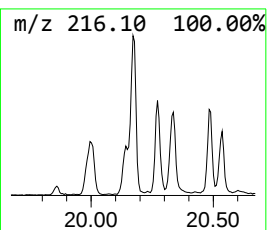
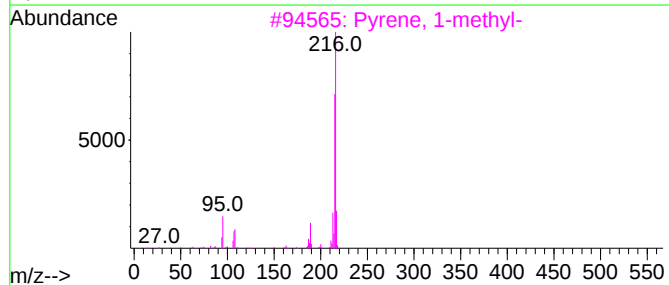
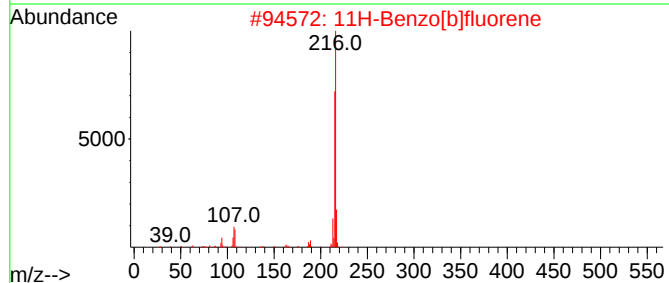
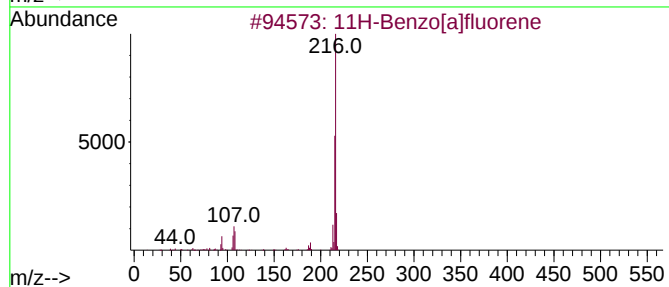
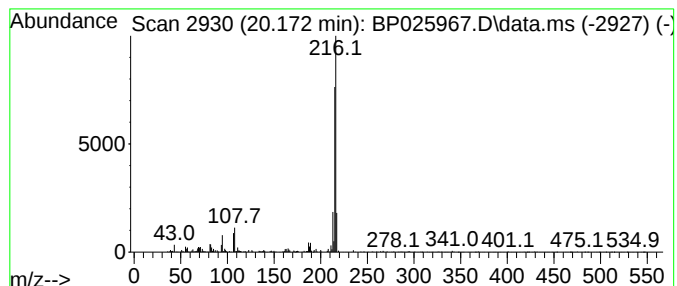
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 11H-Benzo[a]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.172	2.13 ng	400848	Chrysene-d12	21.583

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	97
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	96
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	86
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100925\
Data File : BP025967.D
Acq On : 09 Oct 2025 16:27
Operator : RC/JU
Sample : Q3313-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETGI-360

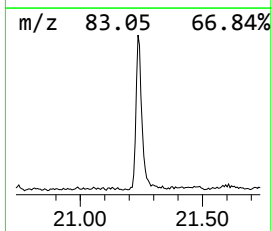
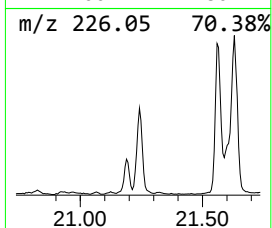
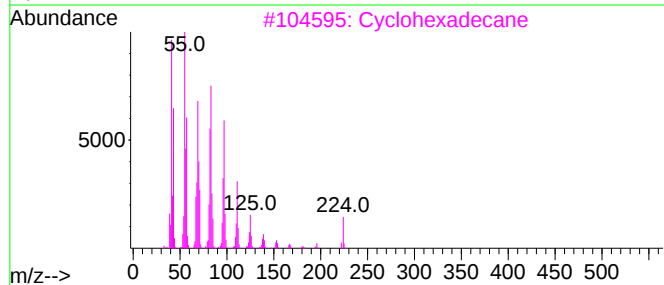
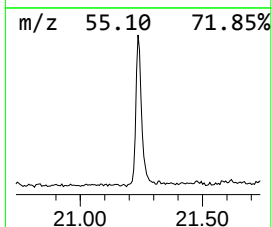
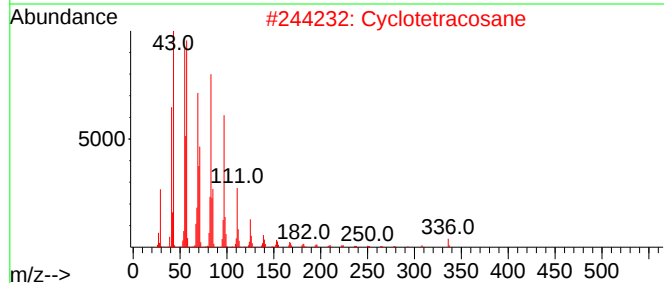
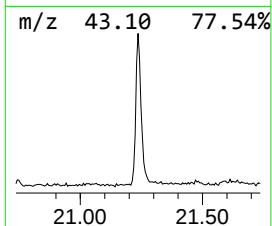
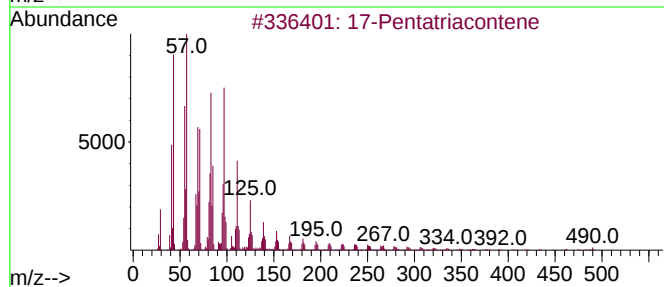
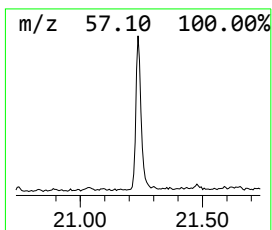
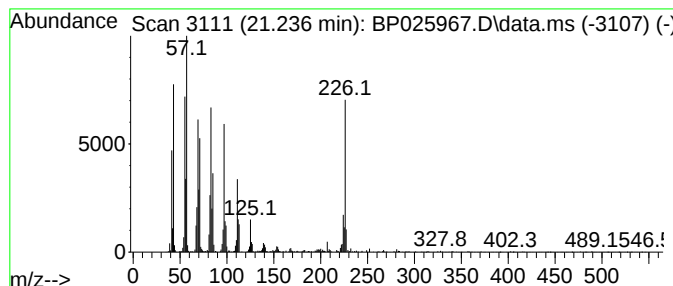
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 17-Pentatriacontene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.236	5.71 ng	1074720	Chrysene-d12	21.583

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			17-Pentatriacontene	491	C35H70	006971-40-0	93
2			Cyclotetracosane	336	C24H48	000297-03-0	92
3			Cyclohexadecane	224	C16H32	000295-65-8	78
4			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	70
5			Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100925\
Data File : BP025967.D
Acq On : 09 Oct 2025 16:27
Operator : RC/JU
Sample : Q3313-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETGI-360

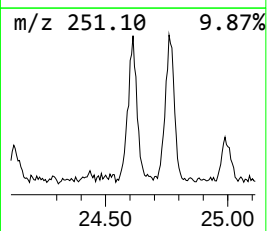
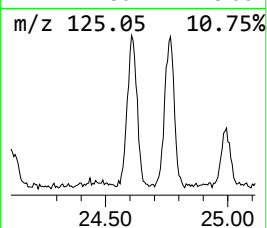
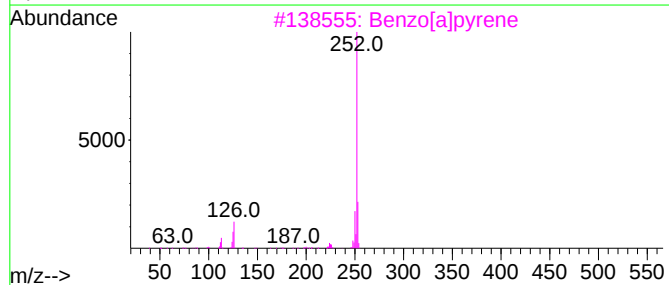
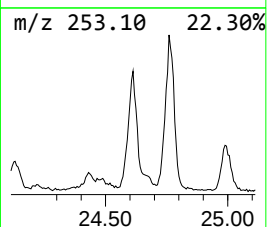
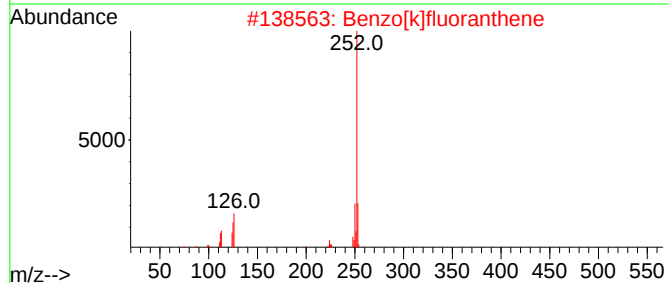
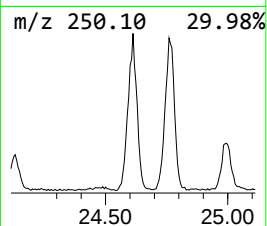
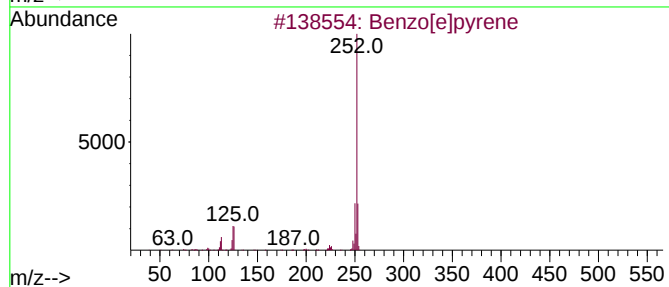
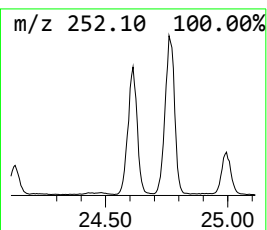
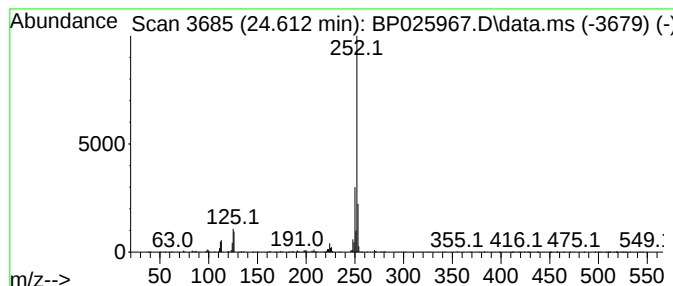
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.613	9.78 ng	958230	Perylene-d12	24.918

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
3			Benzo[a]pyrene	252	C20H12	000050-32-8	96
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	95
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100925\
Data File : BP025967.D
Acq On : 09 Oct 2025 16:27
Operator : RC/JU
Sample : Q3313-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETGI-360

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.713	7.1	ng	637428	3	14.342	1798580	20.0
n-Hexadecanoic ...	18.066	16.8	ng	1748360	4	17.142	2087130	20.0
4H-Cyclopenta[d...]	18.248	3.3	ng	341605	4	17.142	2087130	20.0
Octadecanoic acid	19.401	2.2	ng	418353	5	21.583	3765150	20.0
11H-Benzo[a]flu...	20.172	2.1	ng	400848	5	21.583	3765150	20.0
17-Pentatriacon...	21.236	5.7	ng	1074720	5	21.583	3765150	20.0
Benzo[e]pyrene	24.613	9.8	ng	958230	6	24.918	1959170	20.0