

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100725\
 Data File : BP025944.D
 Acq On : 07 Oct 2025 15:55
 Operator : RC/JU
 Sample : Q3295-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EO-02-10062025

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: BP025944.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.002	7	11	39	rVB	1021363	1497849	29.61%	2.434%
2	5.373	406	414	439	rBV	1406357	2334583	46.16%	3.794%
3	6.937	673	680	708	rVB	1065741	2234144	44.17%	3.631%
4	7.731	807	815	822	rBV	562035	916695	18.12%	1.490%
5	8.872	1002	1009	1023	rBV	685605	1446867	28.61%	2.351%
6	9.437	1100	1105	1119	rVB	30668	71605	1.42%	0.116%
7	10.490	1273	1284	1300	rBV	658900	1297300	25.65%	2.108%
8	11.902	1517	1524	1531	rBV2	28991	52639	1.04%	0.086%
9	12.402	1600	1609	1619	rVB7	25458	70666	1.40%	0.115%
10	12.860	1683	1687	1694	rVB4	33255	55056	1.09%	0.089%
11	12.954	1696	1703	1715	rBV	1914772	3246263	64.18%	5.275%
12	13.149	1732	1736	1742	rVB	69247	107216	2.12%	0.174%
13	14.072	1888	1893	1905	rVB	109556	227517	4.50%	0.370%
14	14.243	1917	1922	1928	rBV	113713	167621	3.31%	0.272%
15	14.349	1931	1940	1948	rBV	1001956	1676822	33.15%	2.725%
16	14.413	1948	1951	1959	rVB	53658	92416	1.83%	0.150%
17	14.760	2006	2010	2013	rBV3	39176	52034	1.03%	0.085%
18	15.213	2083	2087	2097	rVB2	124833	194295	3.84%	0.316%
19	15.419	2118	2122	2133	rVB3	56148	134954	2.67%	0.219%
20	15.737	2170	2176	2182	rBV	295369	466535	9.22%	0.758%
21	15.872	2192	2199	2207	rBV	1742541	2671874	52.83%	4.342%
22	16.095	2233	2237	2246	rBV3	122062	295494	5.84%	0.480%
23	16.425	2284	2293	2303	rVB	295119	787281	15.57%	1.279%
24	16.907	2372	2375	2378	rVB	76681	78104	1.54%	0.127%
25	17.154	2410	2417	2421	rBV2	1415270	2057040	40.67%	3.343%
26	17.195	2421	2424	2430	rVV2	1083276	2080530	41.13%	3.381%
27	17.242	2430	2432	2438	rVV	789498	1206235	23.85%	1.960%
28	17.301	2438	2442	2448	rVB	241073	417475	8.25%	0.678%
29	17.460	2466	2469	2472	rBV4	43706	61247	1.21%	0.100%
30	17.666	2502	2504	2507	rBV2	49317	56877	1.12%	0.092%
31	17.854	2533	2536	2541	rBV2	69886	94036	1.86%	0.153%
32	17.972	2552	2556	2564	rBV9	55649	125733	2.49%	0.204%
33	18.066	2568	2572	2573	rBV	154011	167284	3.31%	0.272%
34	18.095	2573	2577	2585	rVV3	1273025	2244716	44.38%	3.648%
35	18.219	2589	2598	2602	rVV2	358856	1001518	19.80%	1.628%
36	18.395	2623	2628	2632	rBV3	150523	284873	5.63%	0.463%

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Min Area: 3 % of largest Peak

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

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37	18.466	2638	2640	2643	rBV3	73905	89891	1.78%	0.146%
38	18.572	2655	2658	2664	rBV	115334	217656	4.30%	0.354%
39	19.019	2729	2734	2737	rBV2	138158	299311	5.92%	0.486%
40	19.284	2772	2779	2787	rBV	3048932	4671743	92.37%	7.592%
41	19.419	2798	2802	2805	rBV	358337	511457	10.11%	0.831%
42	19.660	2834	2843	2852	rVB	2497013	4116073	81.38%	6.689%
43	19.872	2873	2879	2884	rVB	3601095	5057890	100.00%	8.219%
44	20.019	2897	2904	2906	rBV3	173756	397706	7.86%	0.646%
45	20.054	2907	2910	2917	rVB2	259498	425438	8.41%	0.691%
46	20.160	2923	2928	2933	rBV2	492536	1011400	20.00%	1.644%
47	20.348	2957	2960	2964	rVB2	259605	301839	5.97%	0.491%
48	20.489	2981	2984	2987	rVB	164715	178817	3.54%	0.291%
49	20.536	2988	2992	2996	rBV3	212704	405759	8.02%	0.659%
50	21.248	3109	3113	3120	rVB2	649738	1062146	21.00%	1.726%
51	21.501	3152	3156	3160	rVB	460278	604311	11.95%	0.982%
52	21.578	3164	3169	3170	rBV	1867191	2433657	48.12%	3.955%
53	21.642	3177	3180	3195	rVB	1246054	2694339	53.27%	4.378%
54	23.901	3557	3564	3570	rBV	1121591	2868299	56.71%	4.661%
55	24.789	3709	3715	3728	rVB	902573	2497157	49.37%	4.058%
56	24.960	3738	3744	3751	rVB	686420	1717533	33.96%	2.791%

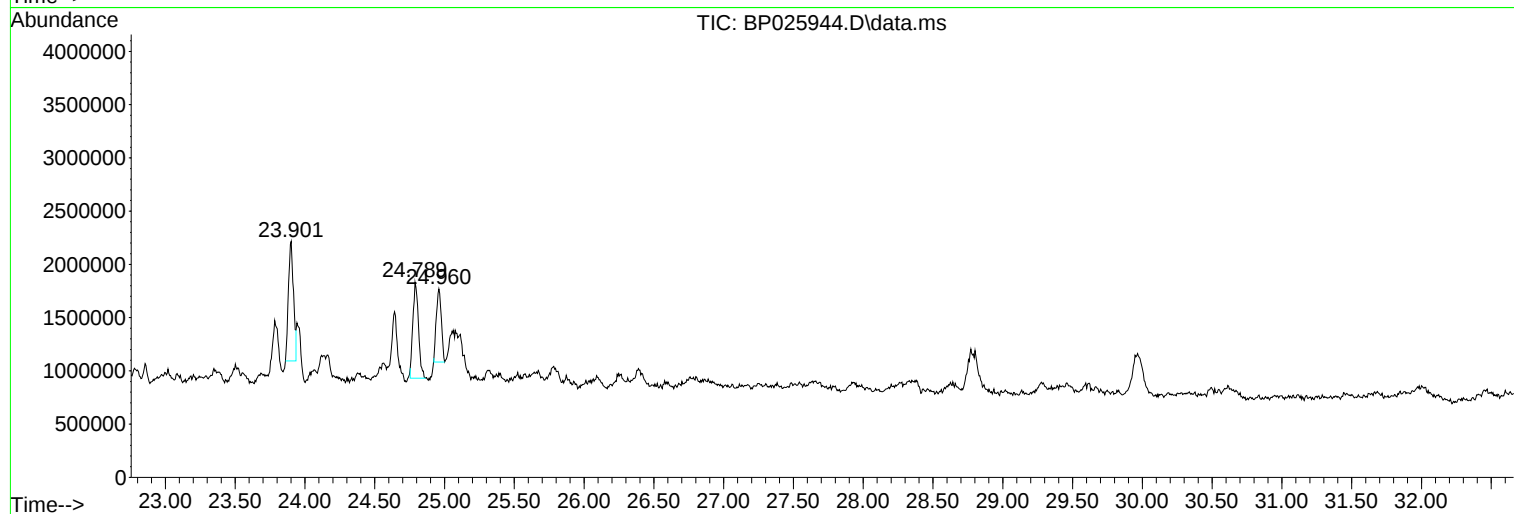
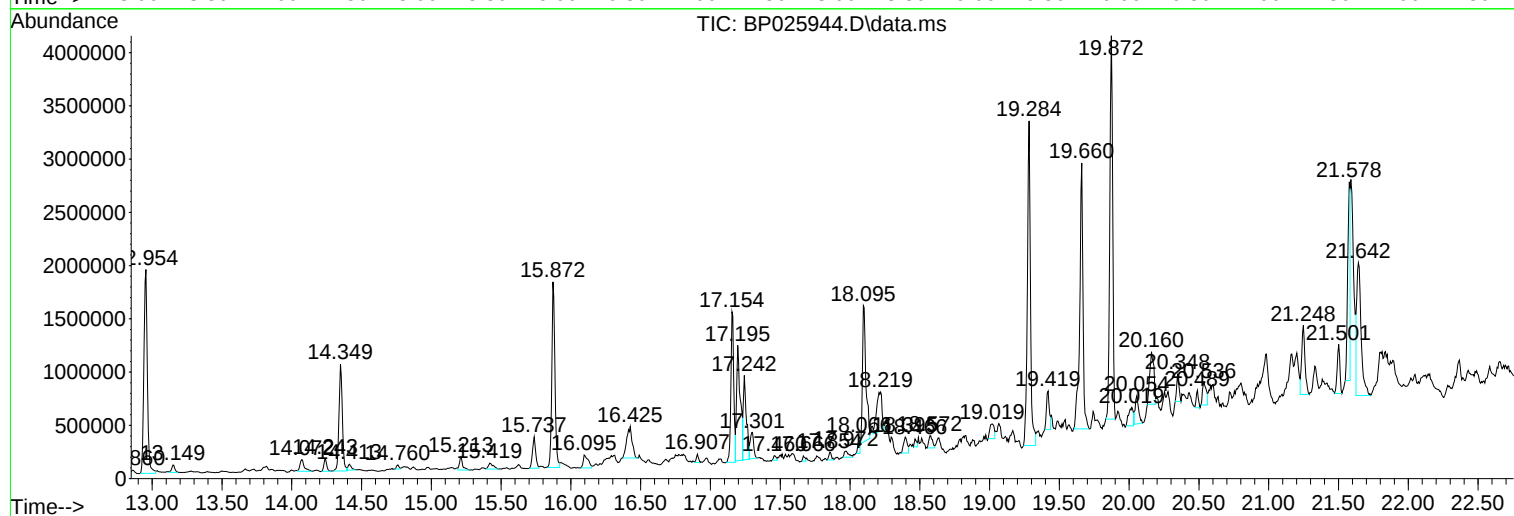
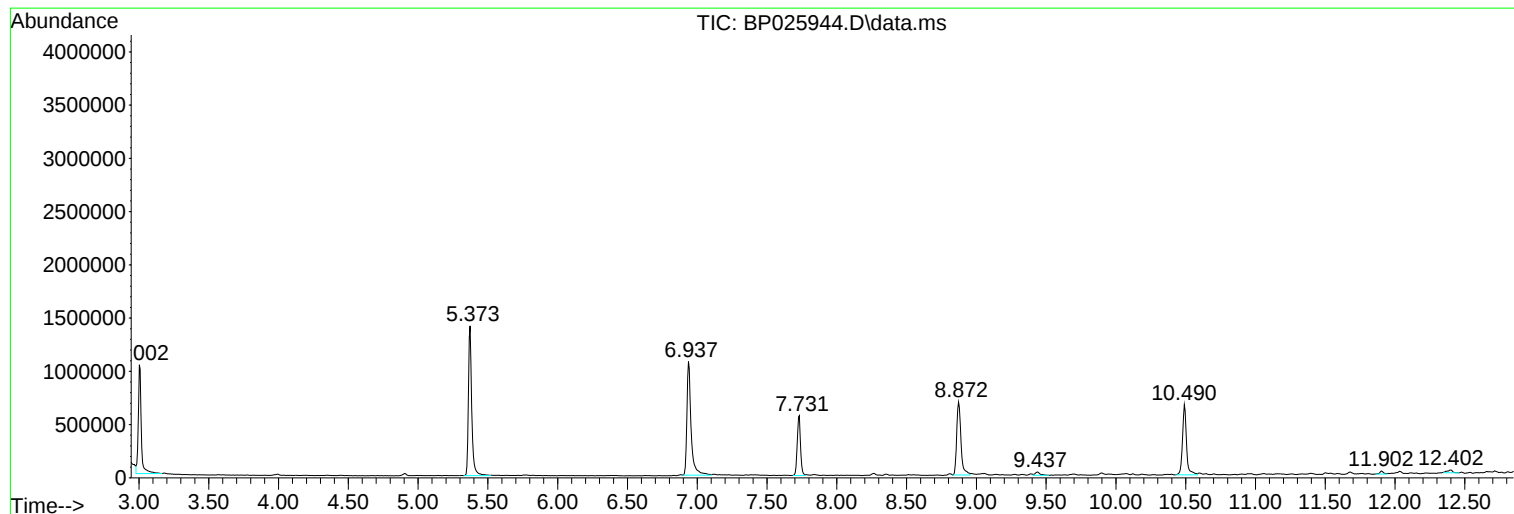
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Operator : RC/JU
Sample : Q3295-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EO-02-10062025

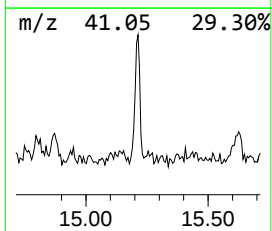
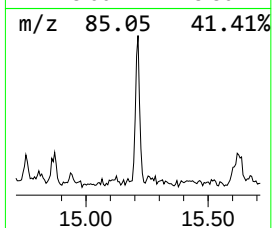
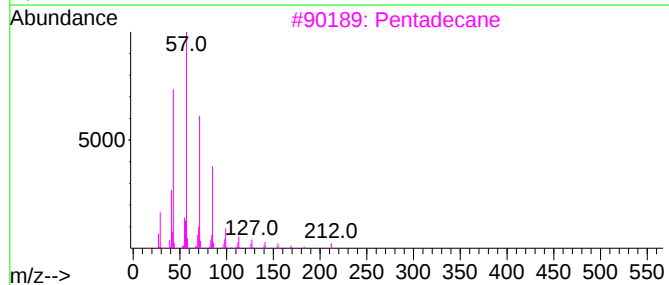
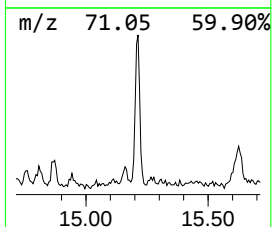
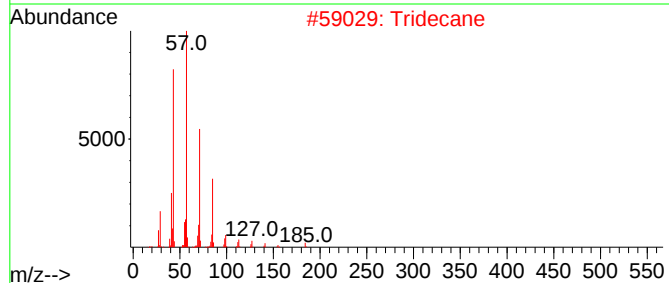
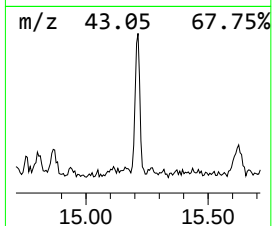
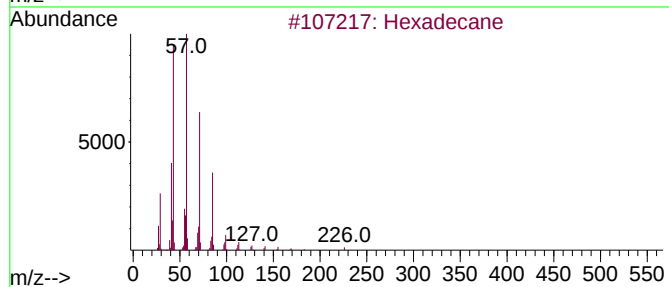
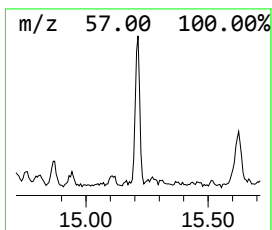
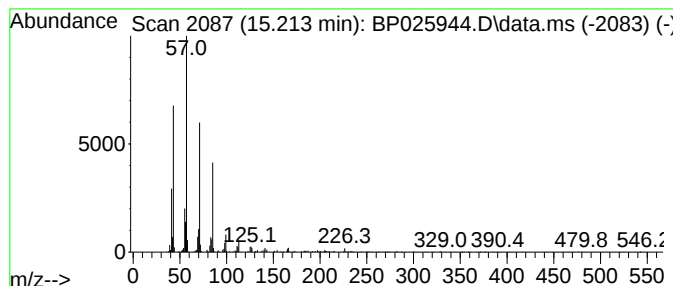
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Hexadecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.213	2.32 ng	194295	Acenaphthene-d10	14.349

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	94
2			Tridecane	184	C13H28	000629-50-5	90
3			Pentadecane	212	C15H32	000629-62-9	90
4			Tetradecane	198	C14H30	000629-59-4	83
5			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	83



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ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EO-02-10062025

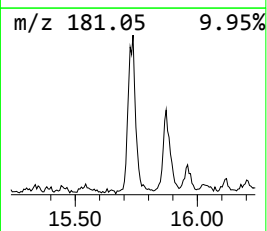
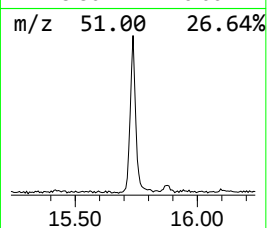
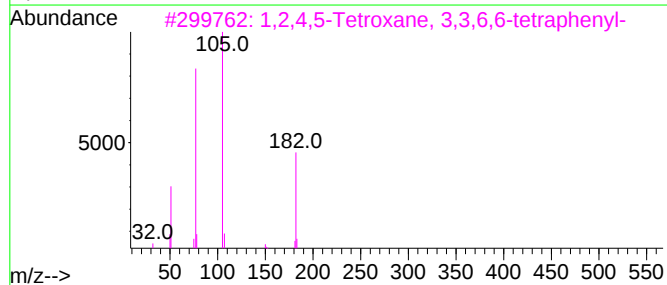
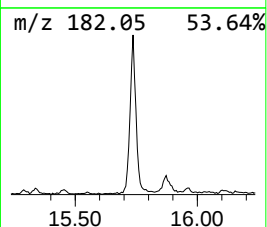
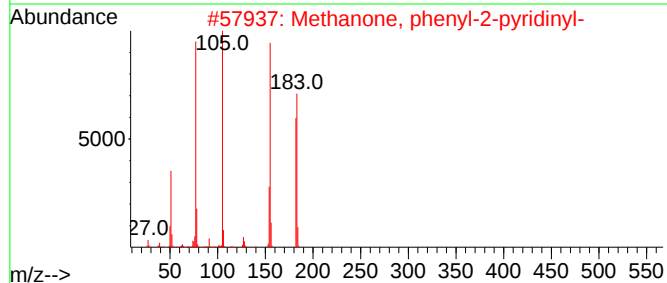
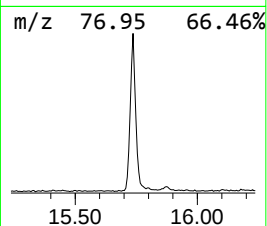
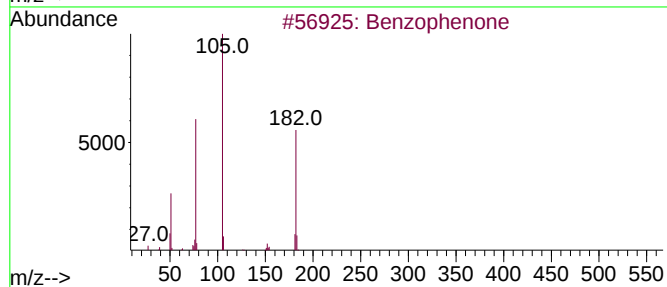
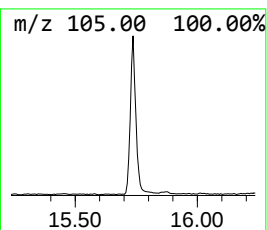
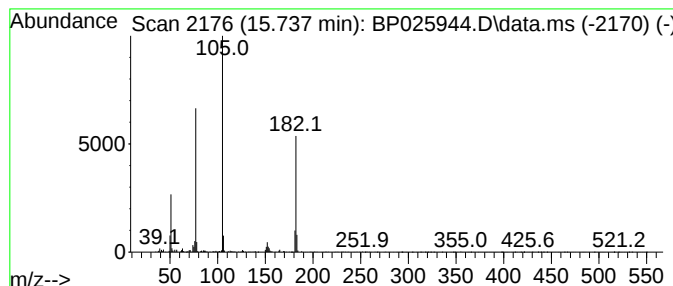
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzophenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.737	5.56 ng	466535	Acenaphthene-d10	14.349

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			1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	64
4			Azobenzene	182	C12H10N2	000103-33-3	37
5			2,3-Dihydro-1-oxo-1H-phenalene	182	C13H10O	000518-85-4	27



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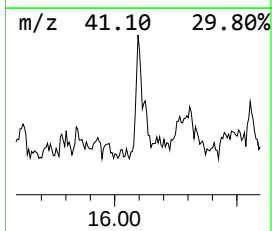
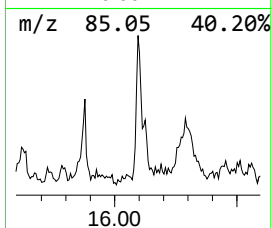
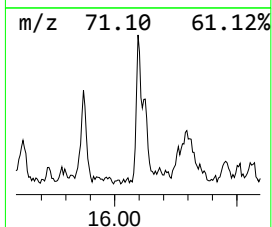
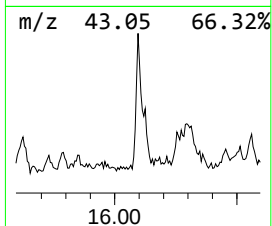
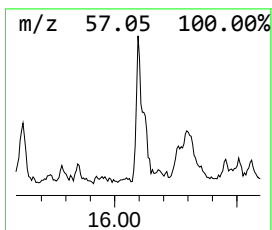
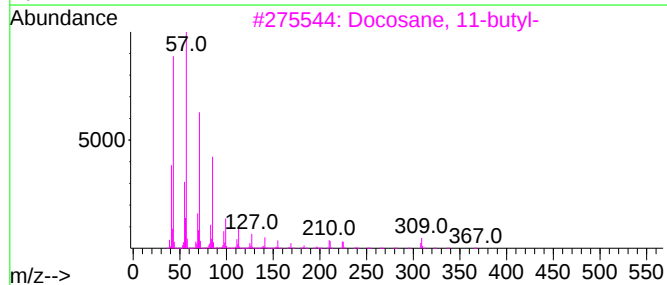
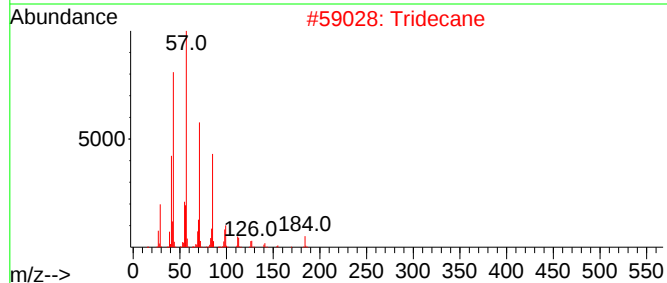
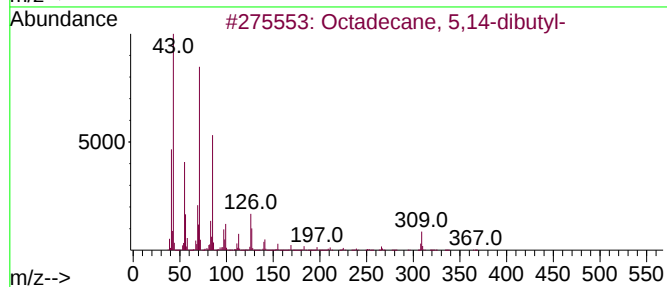
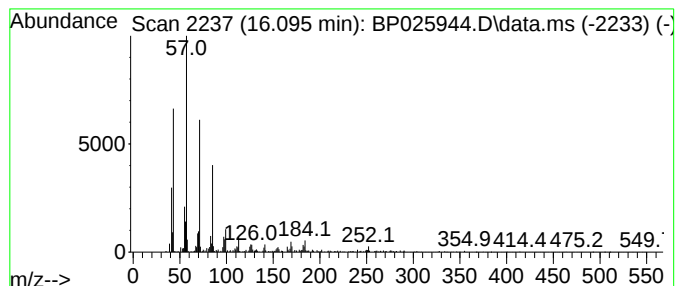
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Octadecane, 5,14-dibutyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.096	2.87 ng	295494	Phenanthrene-d10		17.154	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octadecane, 5,14-dibutyl-		366	C26H54	055282-13-8	93
2	Tridecane		184	C13H28	000629-50-5	93
3	Docosane, 11-butyl-		366	C26H54	013475-76-8	86
4	Tetradecane, 4-ethyl-		226	C16H34	055045-14-2	83
5	Dodecane		170	C12H26	000112-40-3	83



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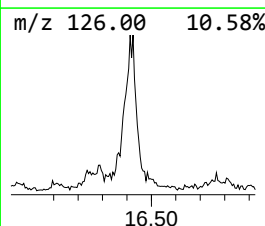
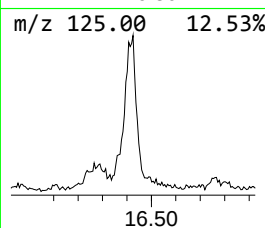
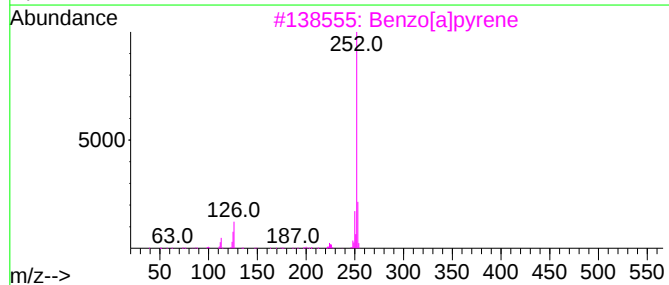
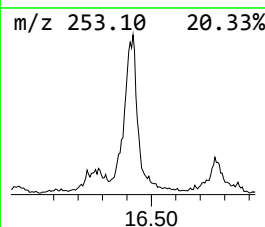
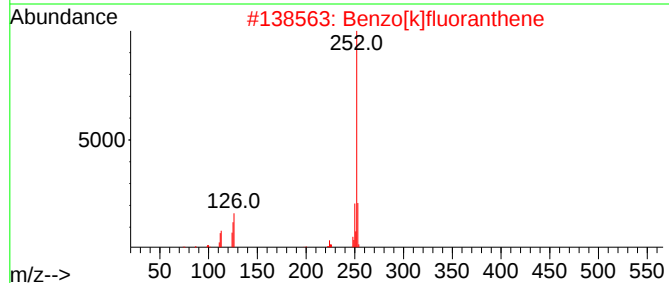
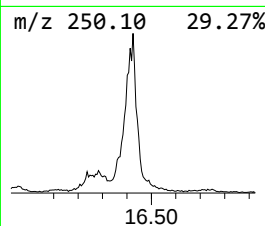
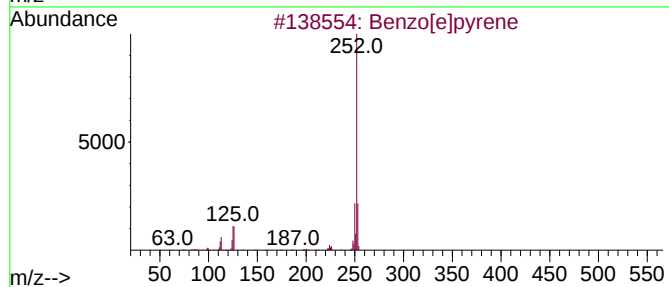
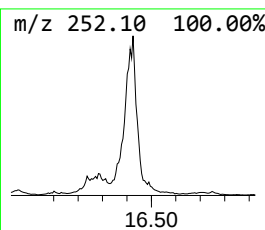
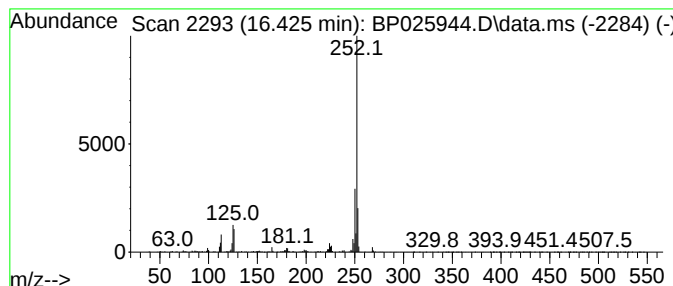
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.425	7.65 ng	787281	Phenanthrene-d10	17.154

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	98
3			Benzo[a]pyrene	252	C20H12	000050-32-8	97
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	90
5			Perylene	252	C20H12	000198-55-0	90



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ALS Vial : 6 Sample Multiplier: 1

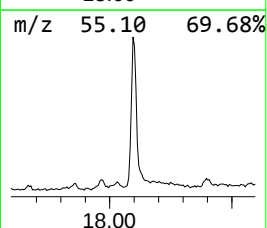
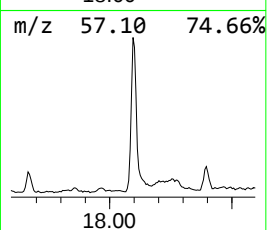
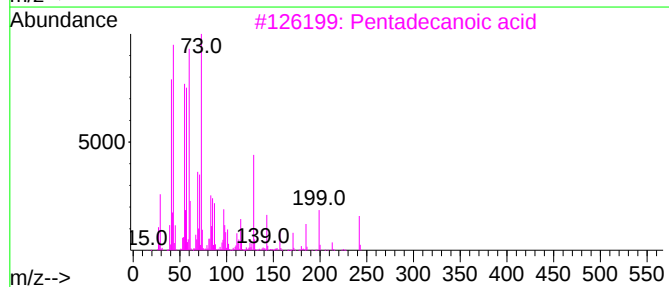
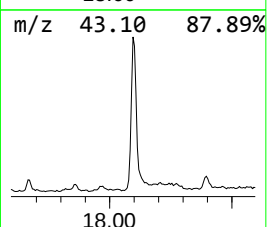
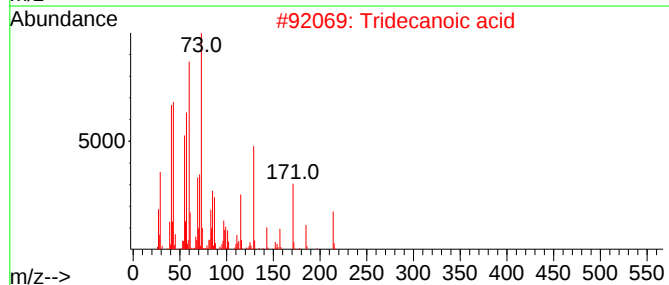
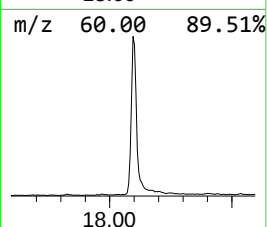
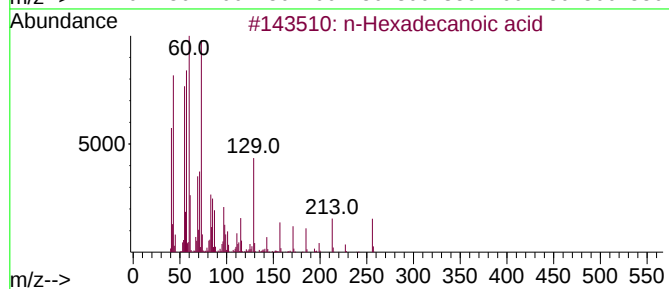
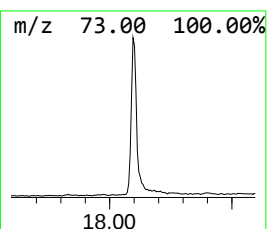
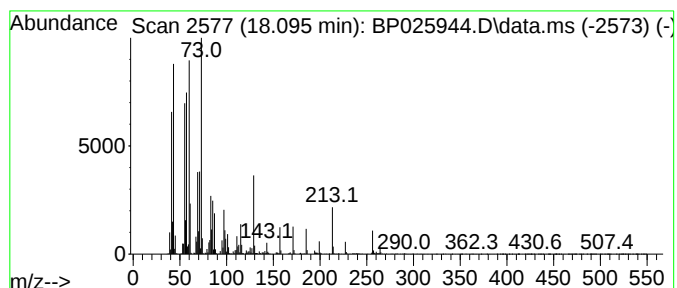
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ClientSampleId :
EO-02-10062025

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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 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.095	21.82 ng	2244720	Phenanthrene-d10	17.154	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	93
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	89
5	Octadecanoic acid	284	C18H36O2	000057-11-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100725\
Data File : BP025944.D
Acq On : 07 Oct 2025 15:55
Operator : RC/JU
Sample : Q3295-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EO-02-10062025

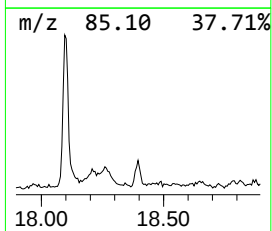
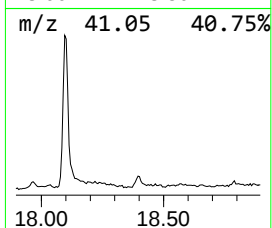
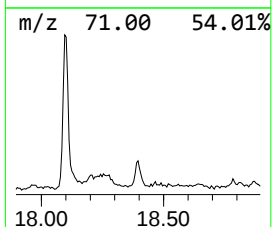
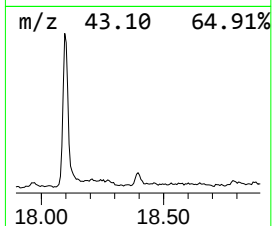
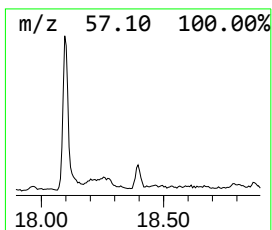
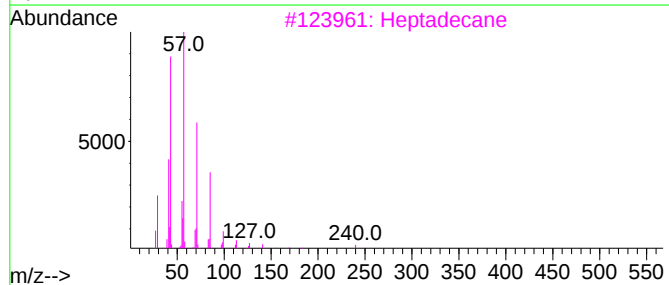
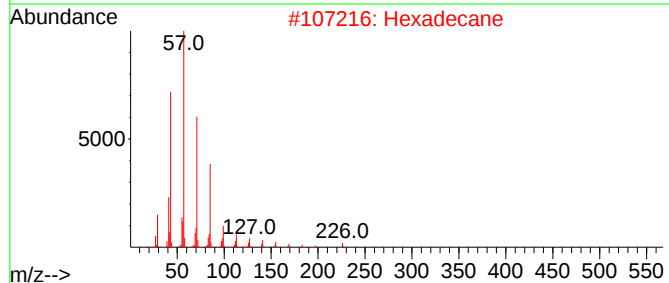
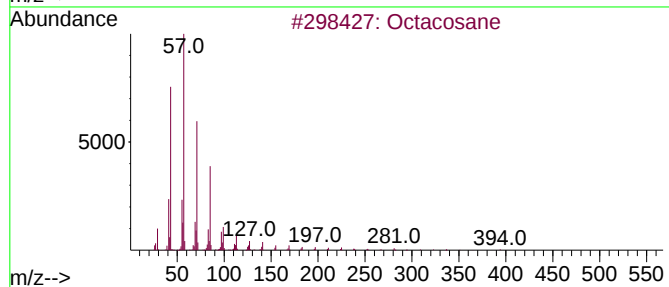
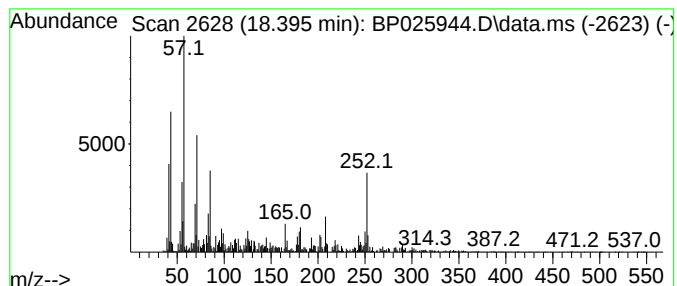
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.395	2.77 ng	284873	Phenanthrene-d10	17.154

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	86
2			Hexadecane	226	C16H34	000544-76-3	83
3			Heptadecane	240	C17H36	000629-78-7	66
4			Tritetracontane	605	C43H88	007098-21-7	41
5			Hexacosane	366	C26H54	000630-01-3	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100725\
Data File : BP025944.D
Acq On : 07 Oct 2025 15:55
Operator : RC/JU
Sample : Q3295-03
Misc :
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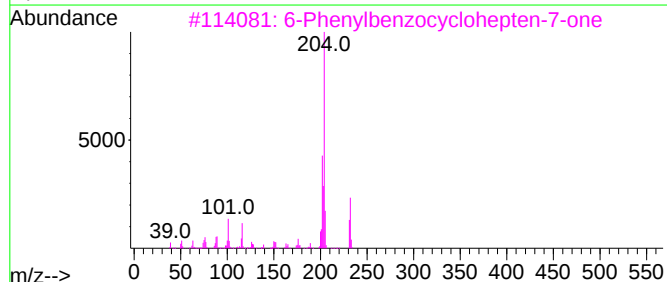
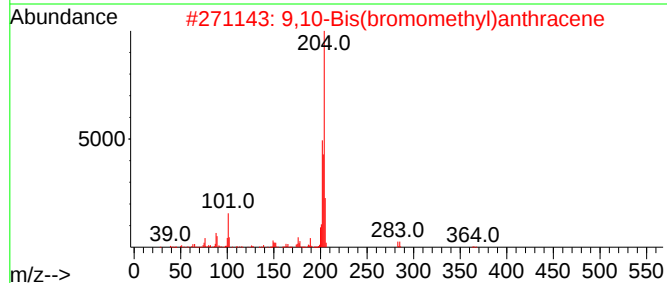
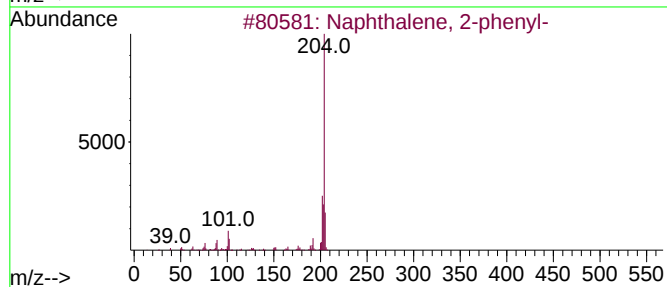
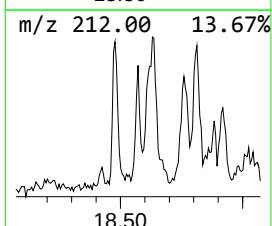
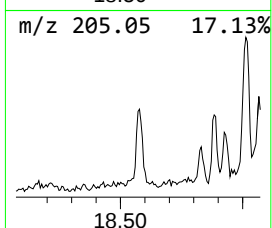
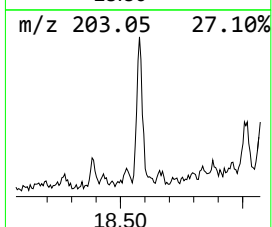
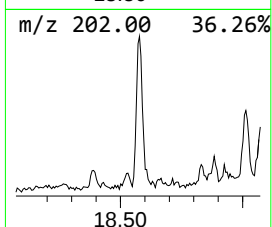
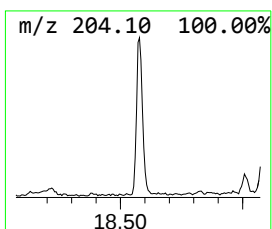
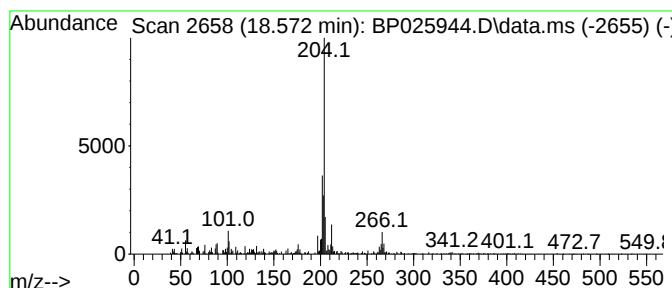
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, 2-phenyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.572	2.12 ng	217656	Phenanthrene-d10	17.154	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	96
2	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	96
3	6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	74
4	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	64
5	5,16[1',2'] :8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100725\
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Acq On : 07 Oct 2025 15:55
Operator : RC/JU
Sample : Q3295-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

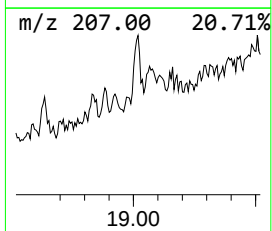
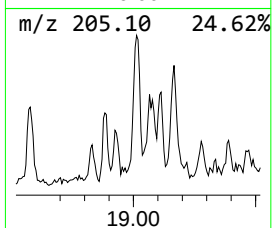
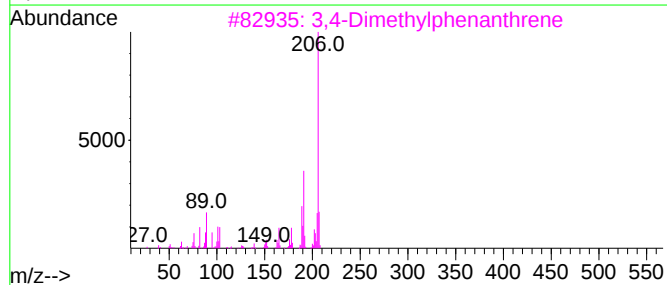
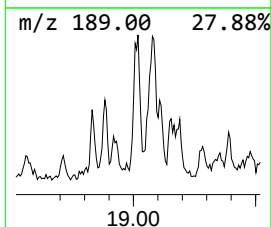
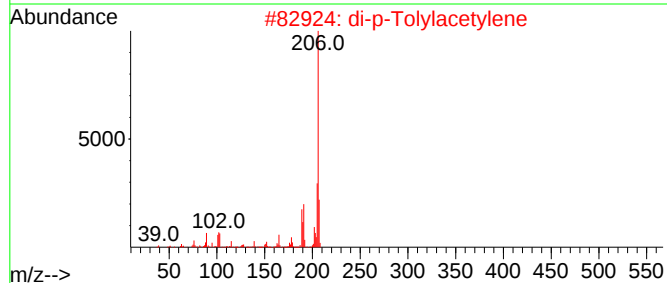
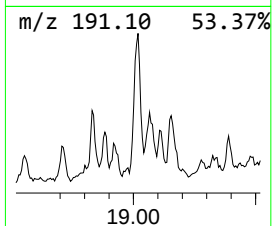
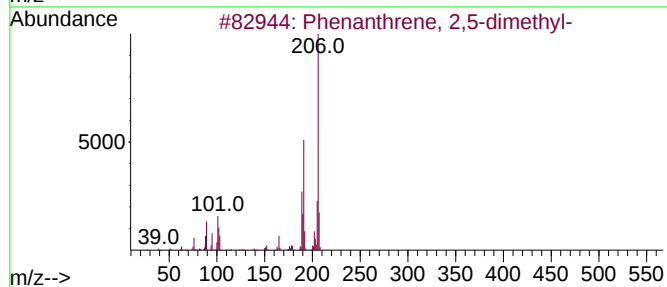
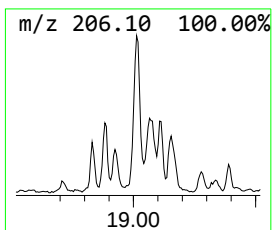
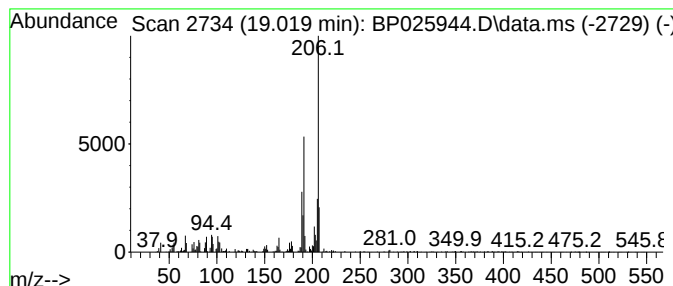
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenanthrene, 2,5-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.019	2.91 ng	299311	Phenanthrene-d10		17.154	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-		206	C16H14	003674-66-6	95
2	di-p-Tolylacetylene		206	C16H14	002789-88-0	93
3	3,4-Dimethylphenanthrene		206	C16H14	1000452-24-5	90
4	Phenanthrene, 3,6-dimethyl-		206	C16H14	001576-67-6	90
5	9,10-Dimethylanthracene		206	C16H14	000781-43-1	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100725\
Data File : BP025944.D
Acq On : 07 Oct 2025 15:55
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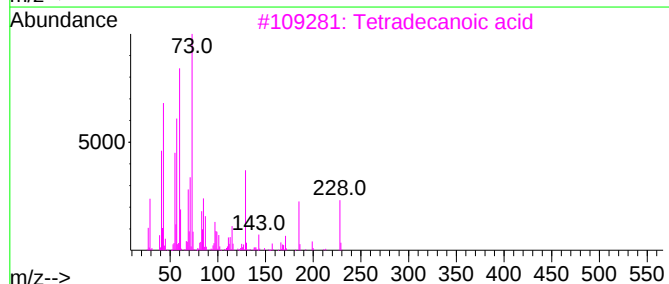
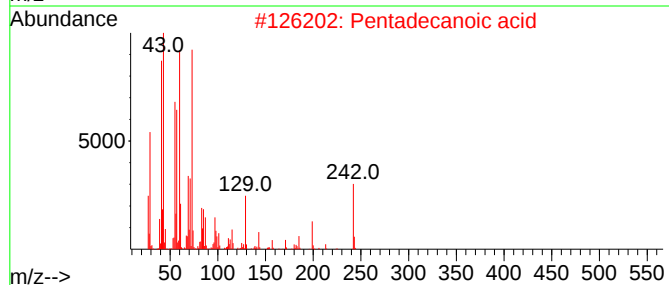
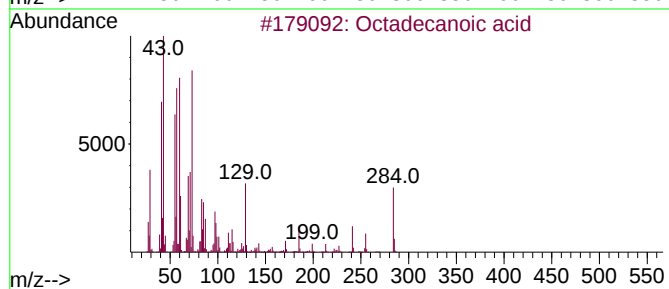
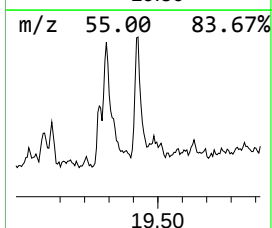
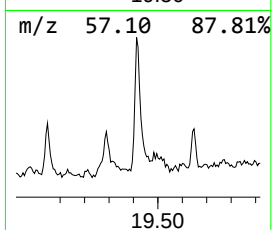
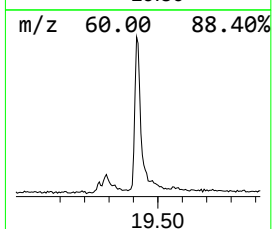
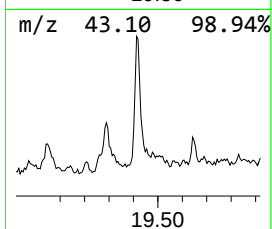
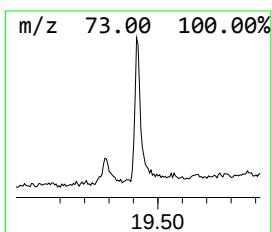
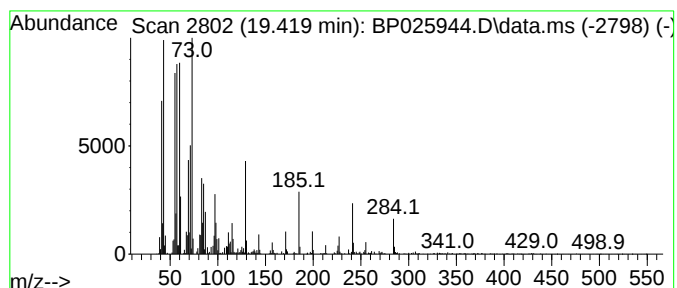
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.419	4.20 ng	511457	Chrysene-d12	21.595

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			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	80
5			n-Decanoic acid	172	C10H20O2	000334-48-5	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100725\
Data File : BP025944.D
Acq On : 07 Oct 2025 15:55
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ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EO-02-10062025

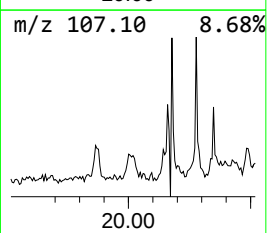
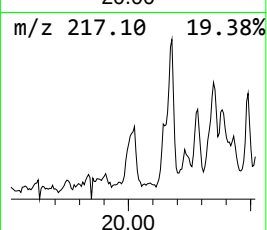
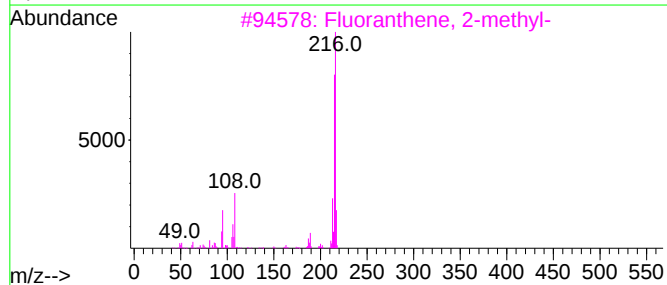
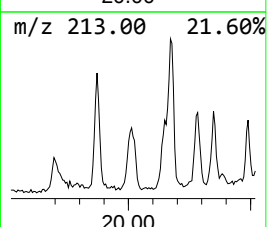
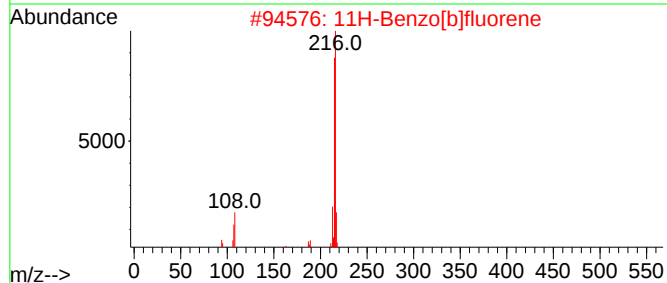
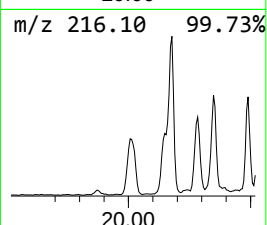
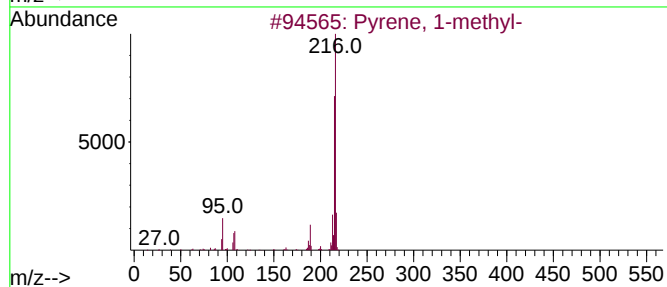
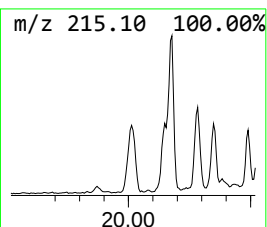
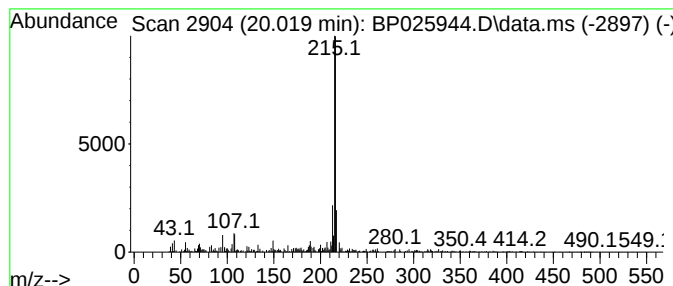
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pyrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.019	3.27 ng	397706	Chrysene-d12	21.595

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
2		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
3		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
4		Pyrene, 2-methyl-	216	C17H12	003442-78-2	68
5		Pyrene, 4-methyl-	216	C17H12	003353-12-6	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100725\
Data File : BP025944.D
Acq On : 07 Oct 2025 15:55
Operator : RC/JU
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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EO-02-10062025

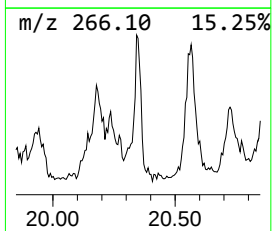
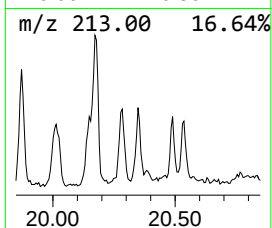
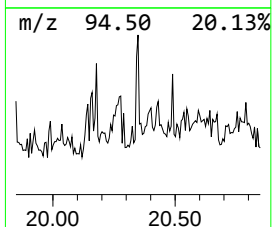
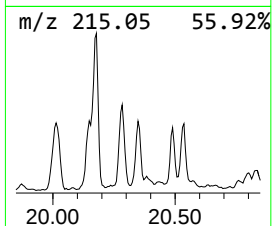
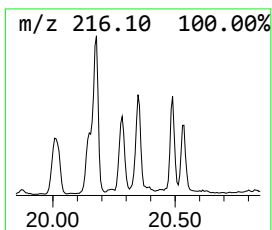
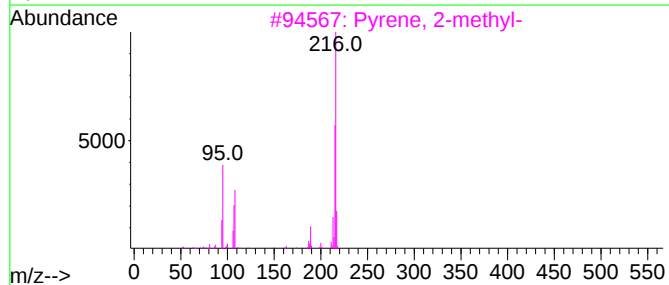
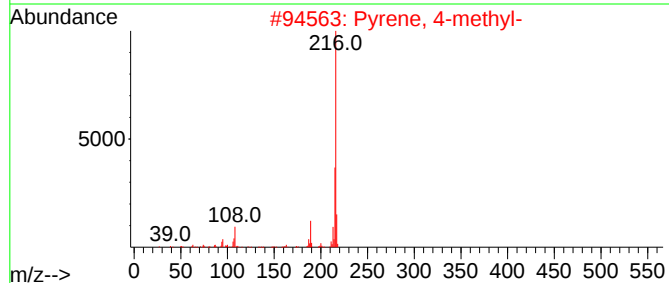
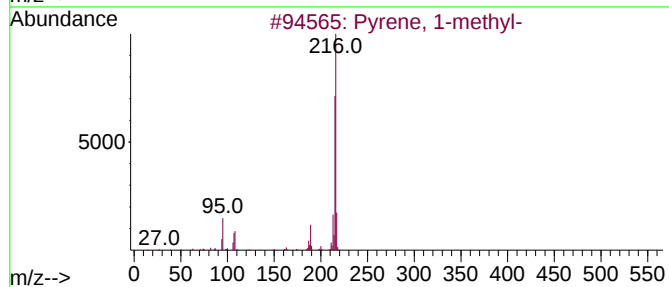
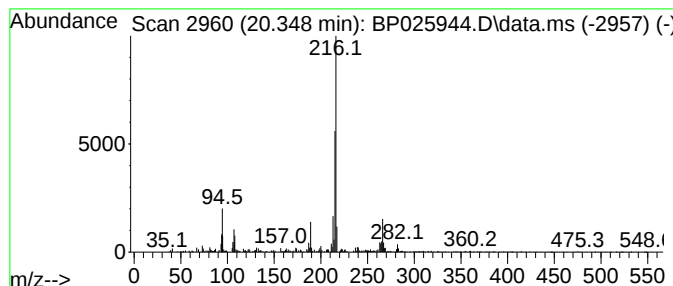
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Pyrene, 4-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.348	2.48 ng	301839	Chrysene-d12	21.595

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	86
2			Pyrene, 4-methyl-	216	C17H12	003353-12-6	60
3			Pyrene, 2-methyl-	216	C17H12	003442-78-2	58
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	58
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100725\
Data File : BP025944.D
Acq On : 07 Oct 2025 15:55
Operator : RC/JU
Sample : Q3295-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EO-02-10062025

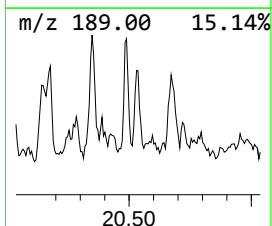
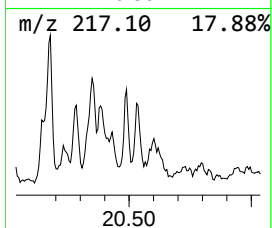
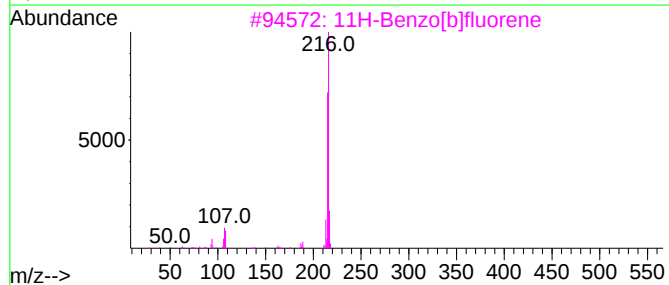
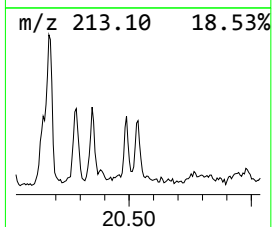
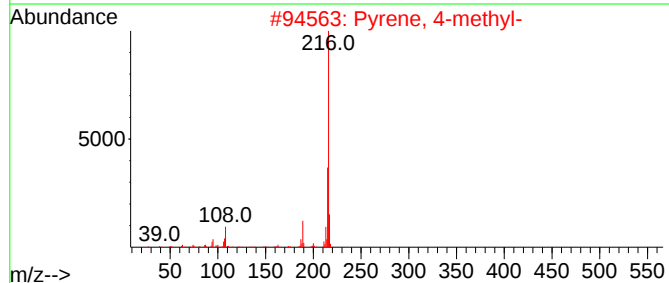
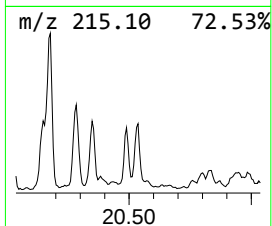
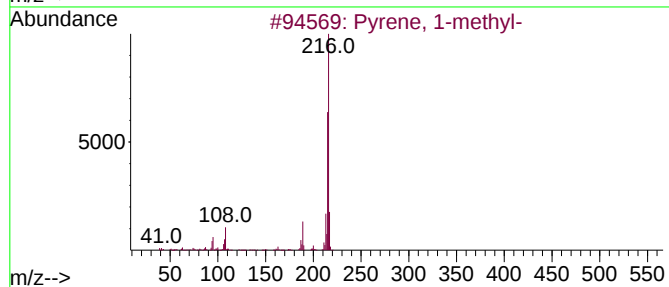
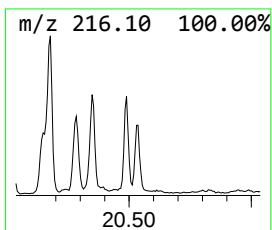
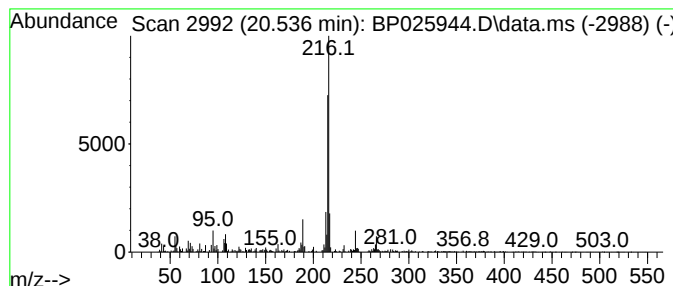
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 11H-Benzo[b]fluorene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.536	3.33 ng	405759	Chrysene-d12	21.595

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-		216	C17H12	002381-21-7	97	
2	Pyrene, 4-methyl-		216	C17H12	003353-12-6	92	
3	11H-Benzo[b]fluorene		216	C17H12	000243-17-4	90	
4	Pyrene, 2-methyl-		216	C17H12	003442-78-2	90	
5	Fluoranthene, 2-methyl-		216	C17H12	033543-31-6	81	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100725\
Data File : BP025944.D
Acq On : 07 Oct 2025 15:55
Operator : RC/JU
Sample : Q3295-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EO-02-10062025

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.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
Hexadecane	15.213	2.3	ng	194295	3	14.349	1676820	20.0
Benzophenone	15.737	5.6	ng	466535	3	14.349	1676820	20.0
Octadecane, 5,1...	16.096	2.9	ng	295494	4	17.154	2057040	20.0
Benzo[e]pyrene	16.425	7.7	ng	787281	4	17.154	2057040	20.0
n-Hexadecanoic ...	18.095	21.8	ng	2244720	4	17.154	2057040	20.0
Octacosane	18.395	2.8	ng	284873	4	17.154	2057040	20.0
Naphthalene, 2-...	18.572	2.1	ng	217656	4	17.154	2057040	20.0
Phenanthrene, 2...	19.019	2.9	ng	299311	4	17.154	2057040	20.0
Octadecanoic acid	19.419	4.2	ng	511457	5	21.595	2433660	20.0
Pyrene, 1-methyl-	20.019	3.3	ng	397706	5	21.595	2433660	20.0
Pyrene, 4-methyl-	20.348	2.5	ng	301839	5	21.595	2433660	20.0
11H-Benzo[b]flu...	20.536	3.3	ng	405759	5	21.595	2433660	20.0