

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032422\
 Data File : BP009534.D
 Acq On : 24 Mar 2022 17:52
 Operator : CG/JU
 Sample : N2090-16
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-60

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-BP031722.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP009534.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.022	3	6	18	rVB	1620894	2175572	8.45%	2.015%
2	5.375	400	406	430	rBV	2093730	3709509	14.41%	3.435%
3	6.963	669	676	697	rBV	1078630	2288846	8.89%	2.120%
4	7.769	807	813	823	rBV	850217	1547562	6.01%	1.433%
5	8.928	1002	1010	1035	rBV	3030869	5644207	21.92%	5.227%
6	10.563	1280	1288	1307	rBV	1072591	2116659	8.22%	1.960%
7	13.040	1700	1709	1731	rBV	8410864	13792250	53.57%	12.773%
8	14.416	1936	1943	1953	rBV	1762354	2928036	11.37%	2.712%
9	15.916	2191	2198	2217	rBV	6424387	10793720	41.92%	9.996%
10	16.181	2237	2243	2257	rBV	634370	997109	3.87%	0.923%
11	17.010	2379	2384	2395	rVB	745168	1221072	4.74%	1.131%
12	17.180	2407	2413	2425	rBV	2231591	3650705	14.18%	3.381%
13	17.622	2484	2488	2494	rBV2	248053	379390	1.47%	0.351%
14	18.098	2566	2569	2579	rVB2	309352	556408	2.16%	0.515%
15	18.239	2590	2593	2597	rVB2	221838	289949	1.13%	0.269%
16	18.833	2690	2694	2698	rBV2	278972	377657	1.47%	0.350%
17	18.951	2710	2714	2718	rBV	423312	592648	2.30%	0.549%
18	19.639	2828	2831	2836	rVB	235918	326607	1.27%	0.302%
19	19.763	2846	2852	2864	rVB	14876978	19668041	76.39%	18.215%
20	21.298	3108	3113	3118	rBV	3100424	4364162	16.95%	4.042%
21	21.421	3127	3134	3156	rBV	15516281	25745663	100.00%	23.844%
22	23.686	3513	3519	3531	rVB2	2275440	4811686	18.69%	4.456%

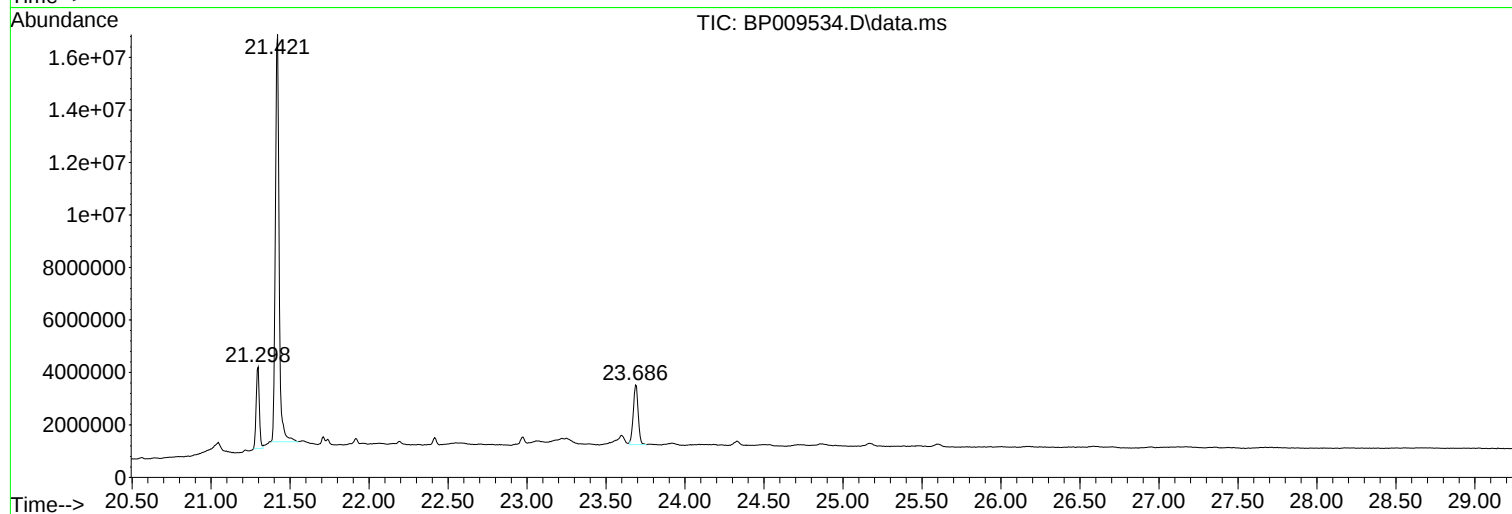
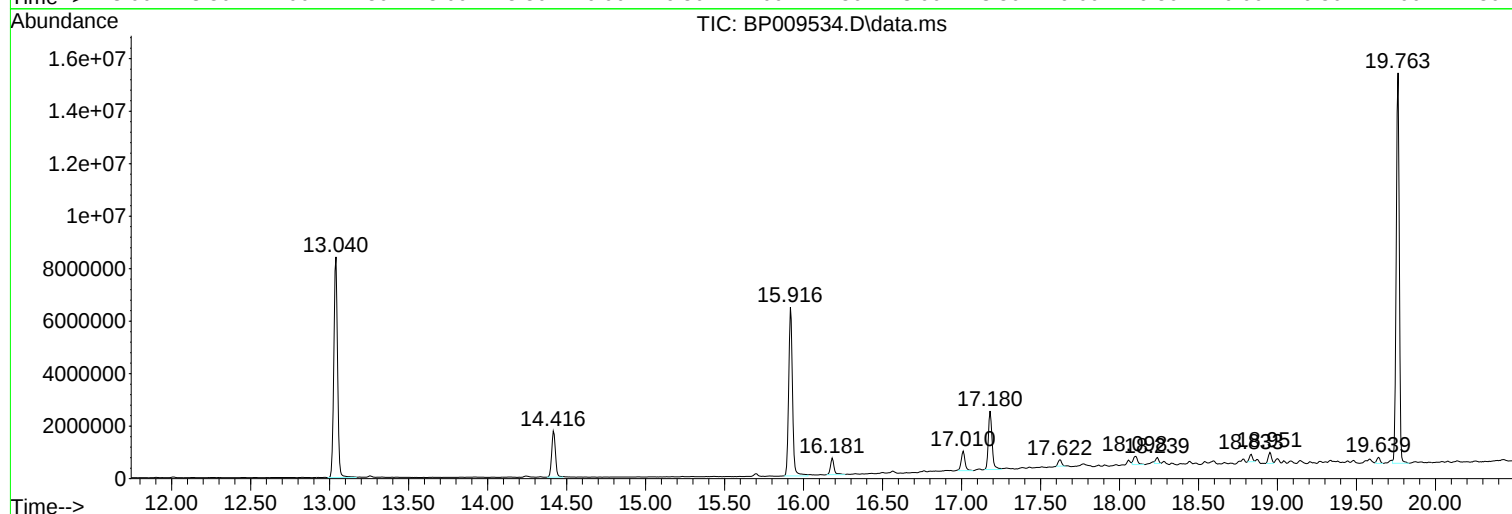
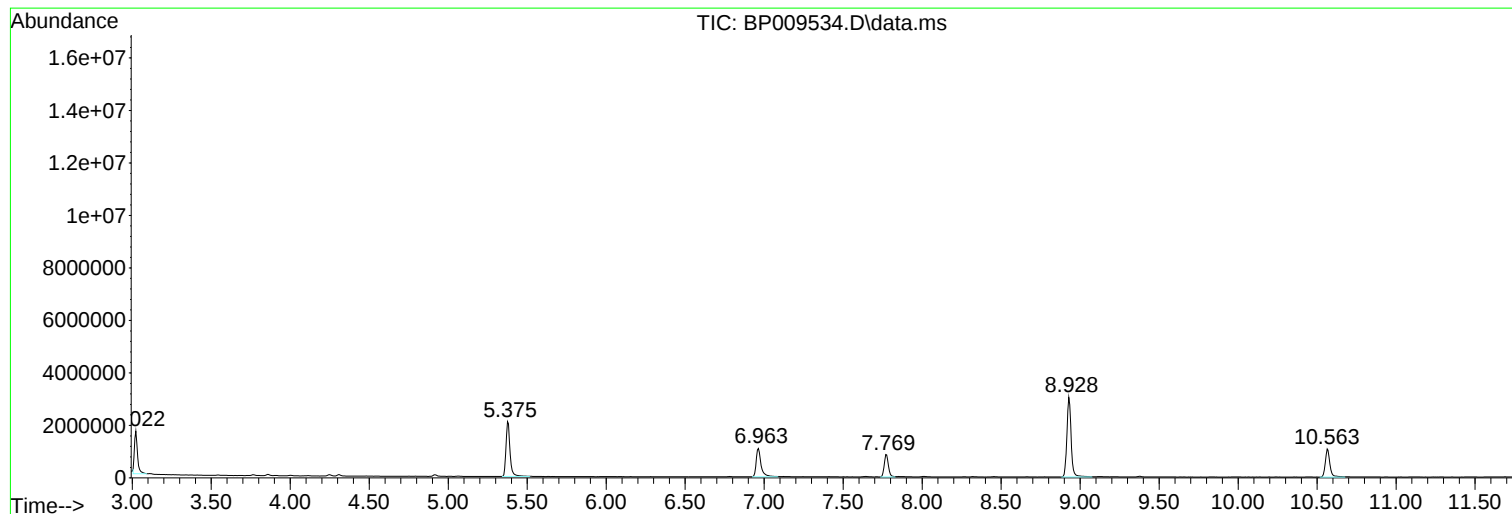
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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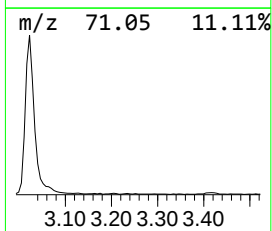
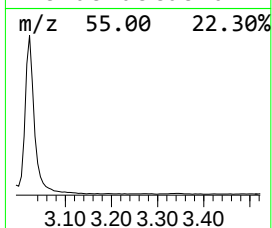
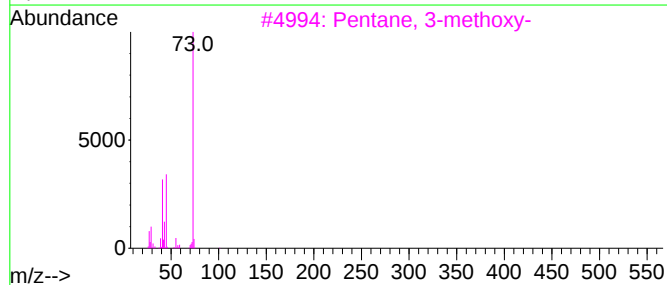
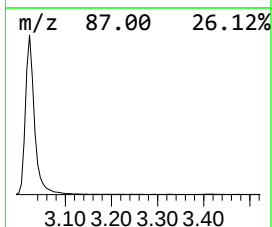
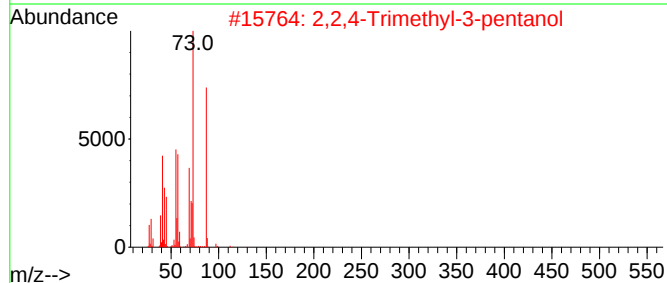
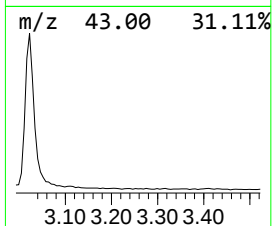
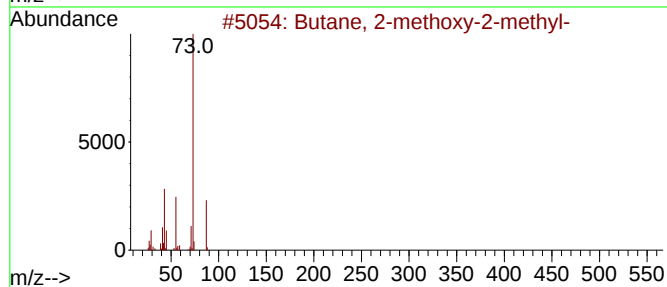
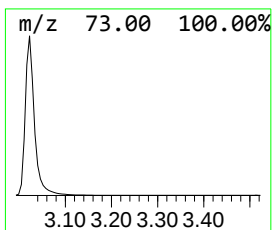
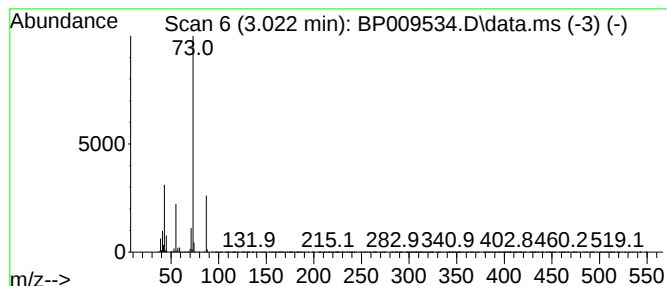
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.022	28.12 ng	2175570	1,4-Dichlorobenzene-d4	7.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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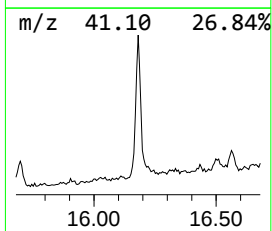
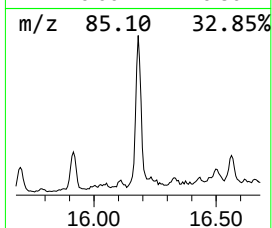
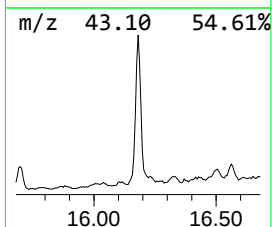
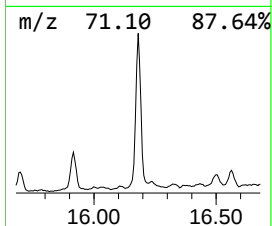
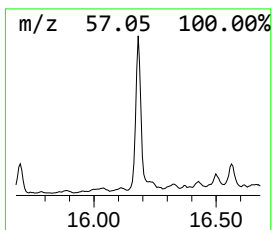
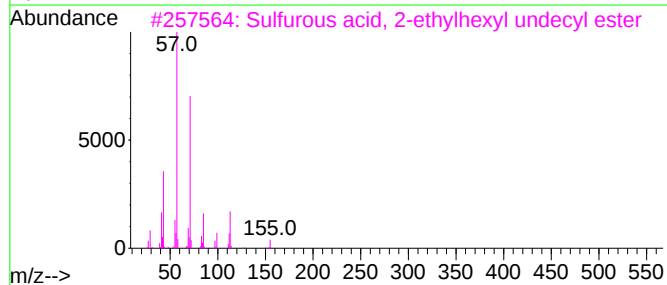
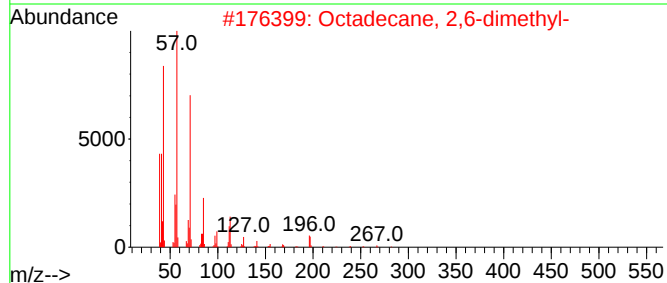
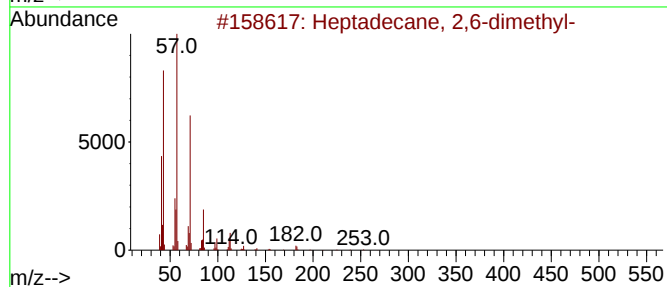
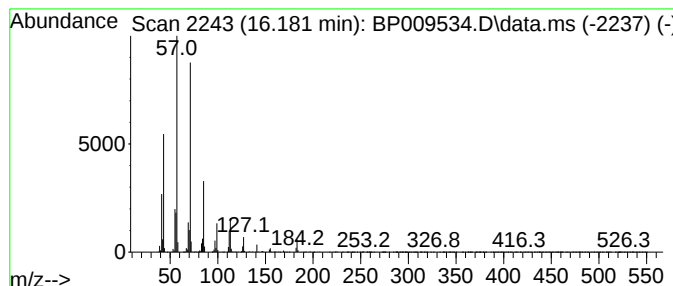
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Heptadecane, 2,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.180	5.46 ng	997109	Phenanthrene-d10	17.180

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	94
2			Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	90
3			Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	80
4			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	74
5			Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	74



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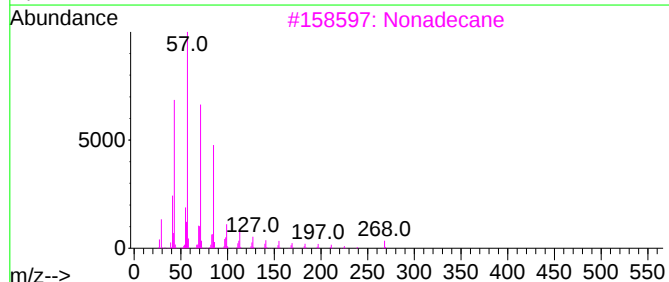
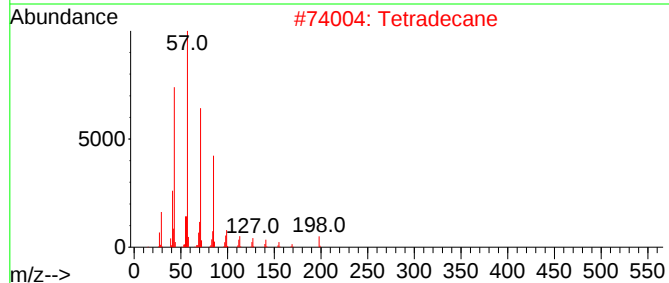
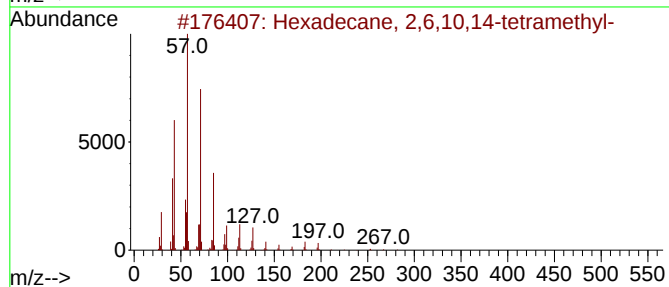
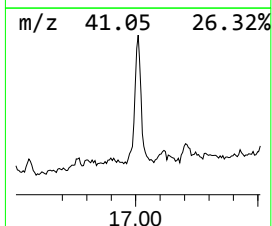
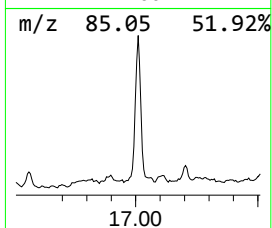
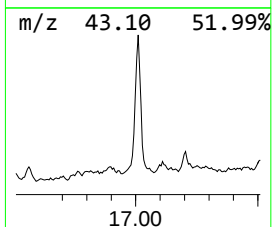
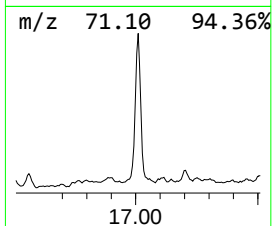
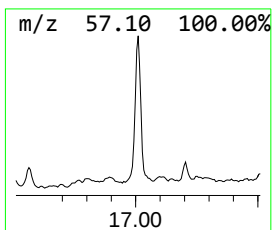
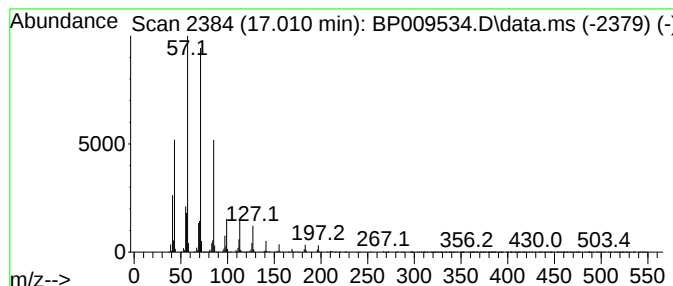
Instrument :
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Hexadecane, 2,6,10,14-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.010	6.69 ng	1221070	Phenanthrene-d10	17.180	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	98
2	Tetradecane	198	C14H30	000629-59-4	90
3	Nonadecane	268	C19H40	000629-92-5	86
4	Pentadecane, 4-methyl-	226	C16H34	002801-87-8	86
5	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	86



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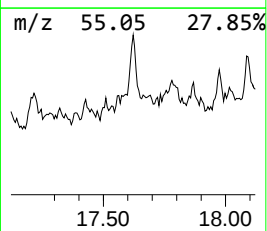
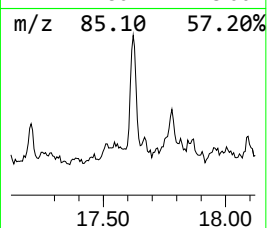
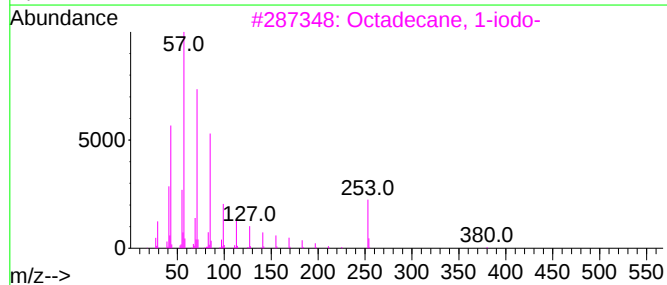
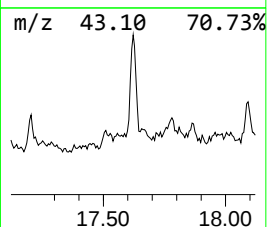
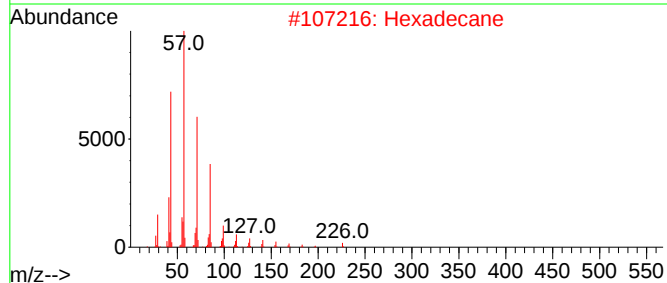
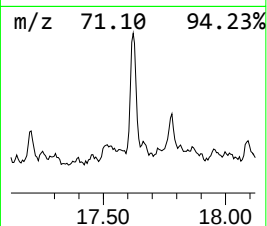
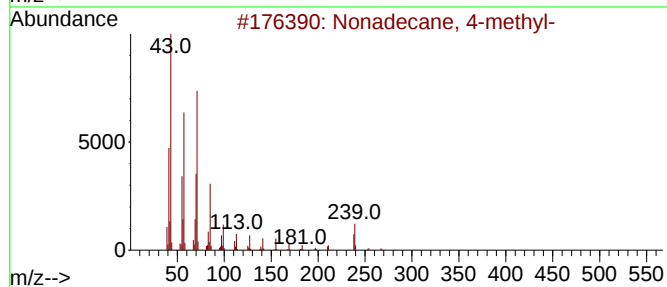
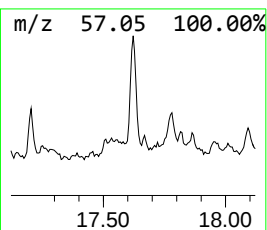
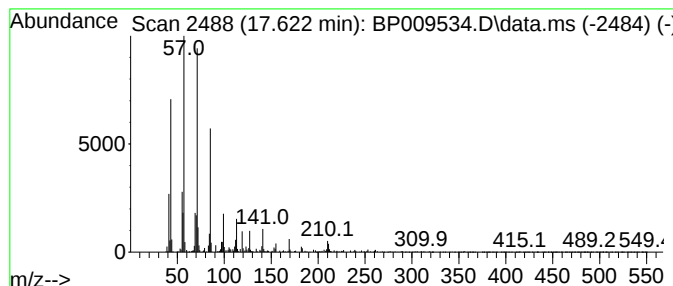
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Peak Number 4 Nonadecane, 4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.622	2.08 ng	379390	Phenanthrene-d10	17.180

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane, 4-methyl-	282	C20H42	025117-27-5	90
2			Hexadecane	226	C16H34	000544-76-3	87
3			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	87
5			2,6,10-Trimethyltridecane	226	C16H34	003891-99-4	87



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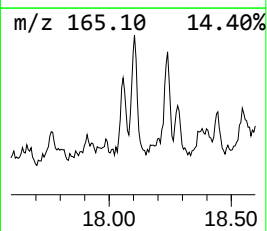
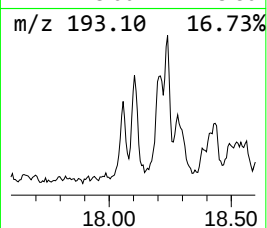
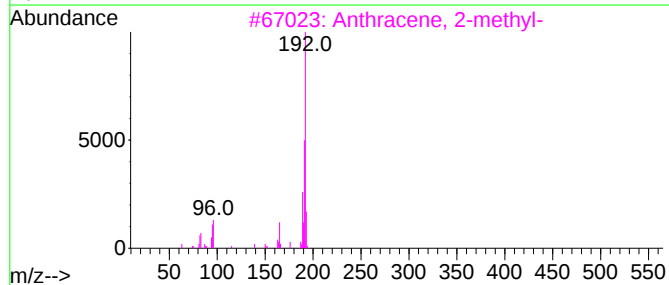
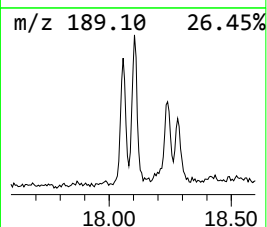
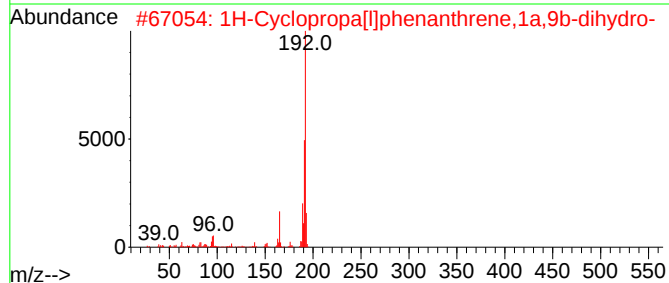
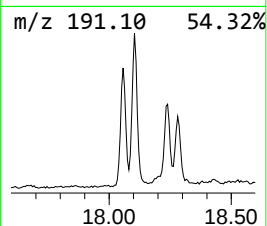
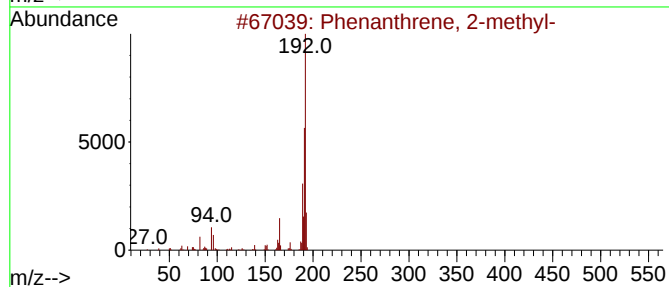
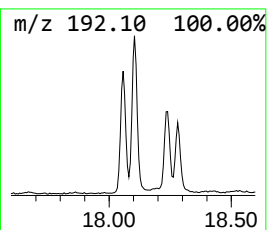
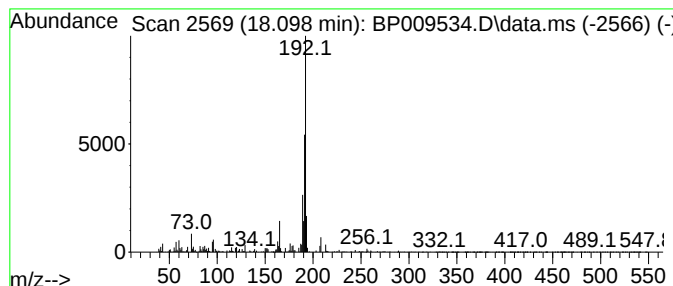
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.098	3.05 ng	556408	Phenanthrene-d10	17.180

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
5			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	90



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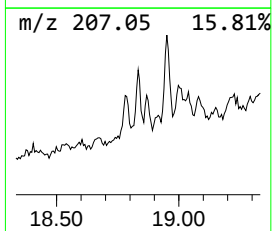
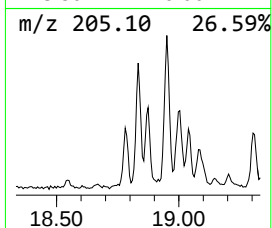
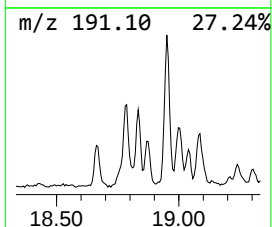
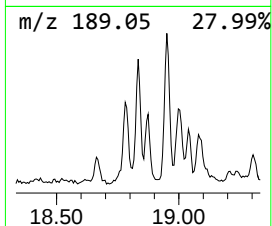
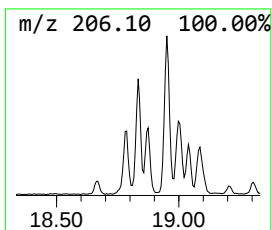
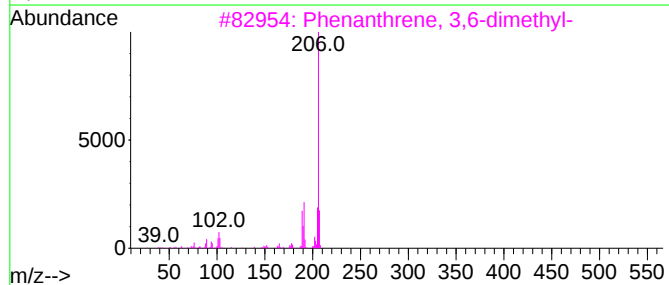
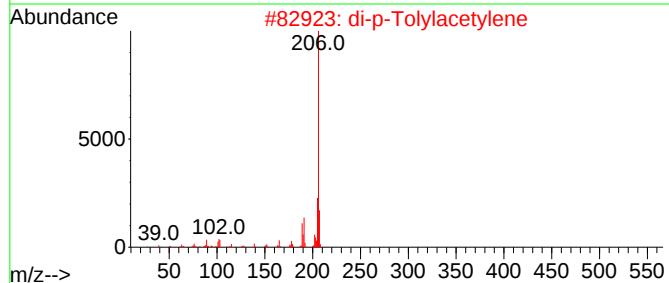
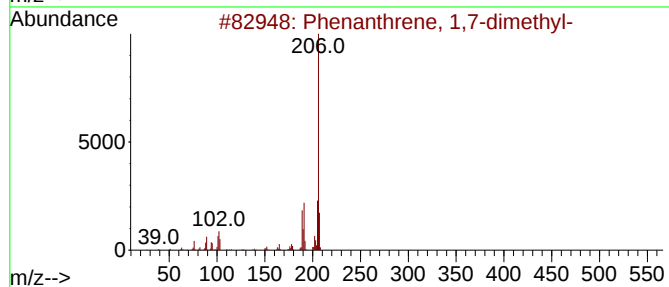
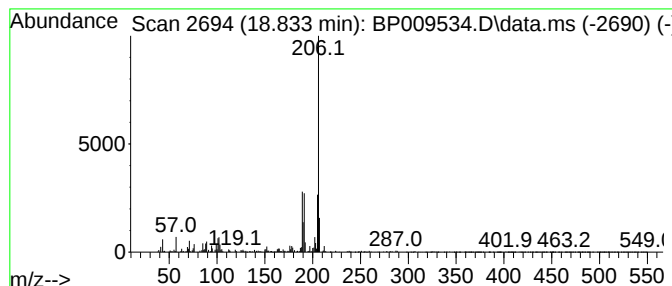
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phenanthrene, 1,7-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.833	2.07 ng	377657	Phenanthrene-d10	17.180

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	94
2			di-p-Tolylacetylene	206	C16H14	002789-88-0	94
3			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
4			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	93
5			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032422\
Data File : BP009534.D
Acq On : 24 Mar 2022 17:52
Operator : CG/JU
Sample : N2090-16
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-60

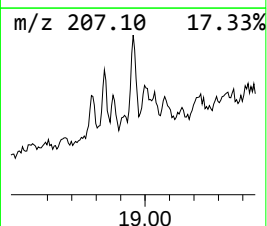
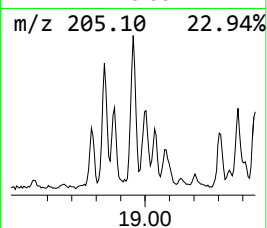
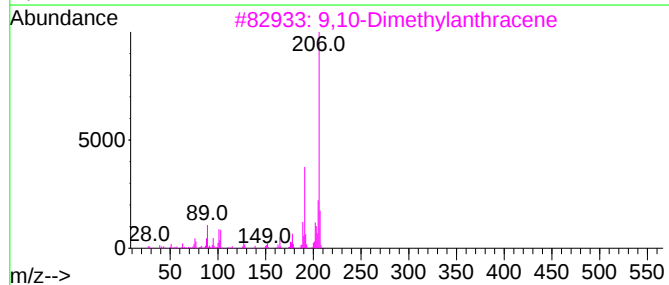
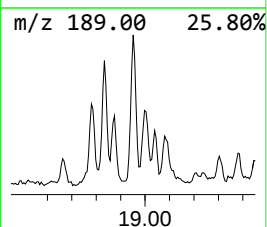
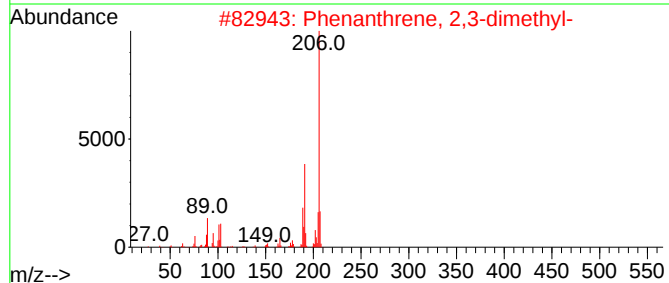
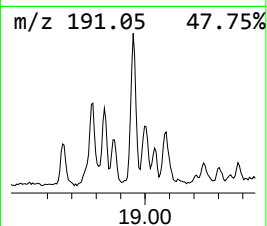
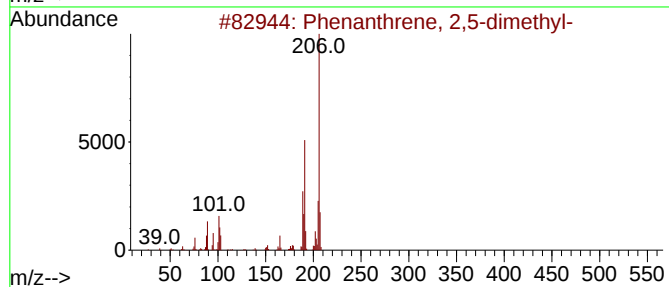
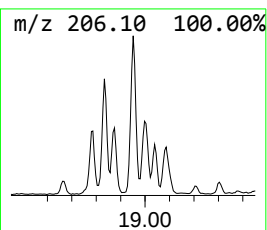
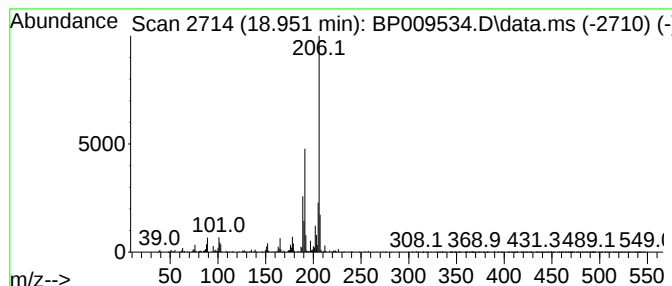
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 2,5-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.951	3.25 ng	592648	Phenanthrene-d10	17.180

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
2			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	93
3			9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
4			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	90
5			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032422\
Data File : BP009534.D
Acq On : 24 Mar 2022 17:52
Operator : CG/JU
Sample : N2090-16
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-60

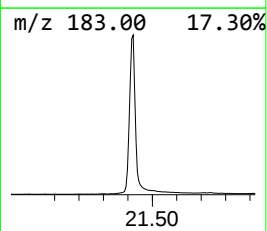
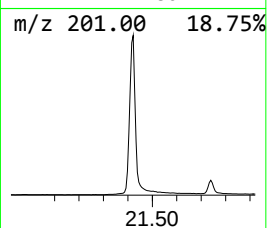
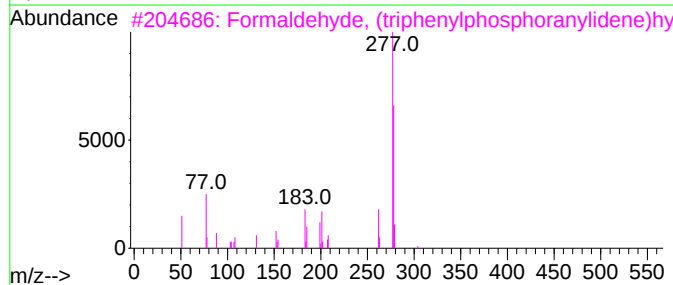
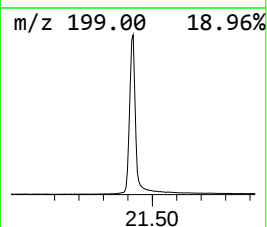
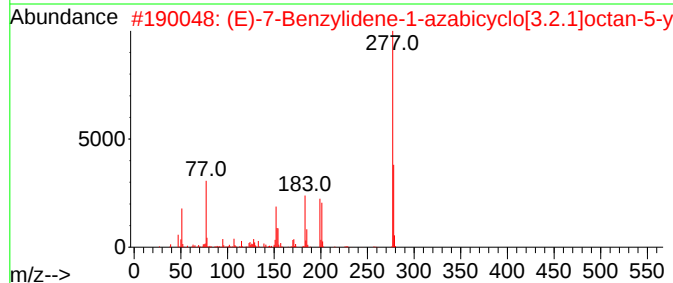
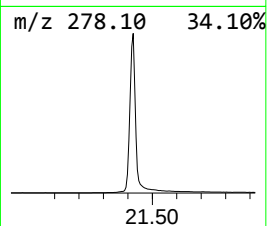
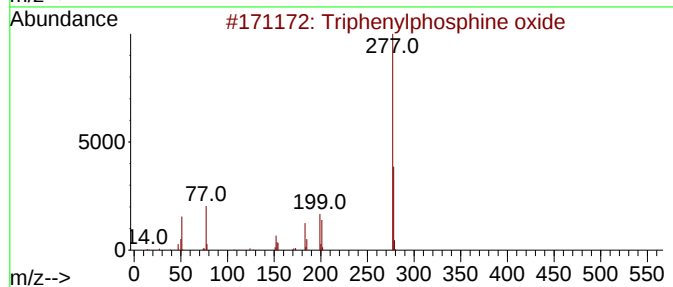
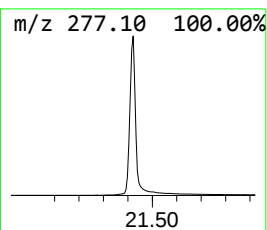
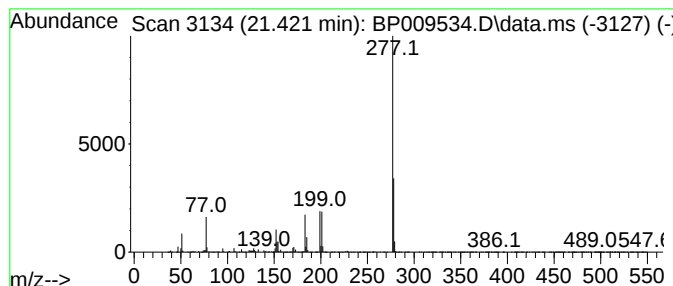
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Triphenylphosphine oxide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.421	117.99 ng	25745700	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	94
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19NO3S	1000483-83-4	91
3			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	47
4			5-Methyl-8-phenylfuro[3',2':5,6]...	277	C16H11N3O2	1000460-07-9	28
5			Thiophene-3-carbonitrile, 4-(4-m...	277	C13H11NO2S2	1000268-75-0	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032422\
Data File : BP009534.D
Acq On : 24 Mar 2022 17:52
Operator : CG/JU
Sample : N2090-16
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-60

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.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
Butane, 2-metho...	3.022	28.1	ng	2175570	1	7.769	1547560	20.0
Heptadecane, 2,...	16.180	5.5	ng	997109	4	17.180	3650710	20.0
Hexadecane, 2,6...	17.010	6.7	ng	1221070	4	17.180	3650710	20.0
Nonadecane, 4-m...	17.622	2.1	ng	379390	4	17.180	3650710	20.0
Phenanthrene, 2...	18.098	3.0	ng	556408	4	17.180	3650710	20.0
Phenanthrene, 1...	18.833	2.1	ng	377657	4	17.180	3650710	20.0
Phenanthrene, 2...	18.951	3.3	ng	592648	4	17.180	3650710	20.0
Triphenylphosph...	21.421	118.0	ng	25745700	5	21.298	4364160	20.0