

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011421\
 Data File : BP004564.D
 Acq On : 14 Jan 2021 19:11
 Operator : CG/JU
 Sample : M1125-01 10X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 1717-1718

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-BP011221.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP004564.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.816	778	787	793	rBV	670770	1103885	31.98%	3.634%
2	7.887	793	799	819	rVB	80586	210799	6.11%	0.694%
3	10.610	1251	1262	1274	rBV	864792	1513591	43.84%	4.983%
4	13.075	1677	1681	1690	rVB	53042	89492	2.59%	0.295%
5	13.810	1802	1806	1811	rBV	108951	168619	4.88%	0.555%
6	13.946	1822	1829	1833	rBV	138624	263685	7.64%	0.868%
7	14.210	1869	1874	1875	rBV4	60356	93689	2.71%	0.308%
8	14.298	1884	1889	1895	rVB4	159804	310438	8.99%	1.022%
9	14.369	1897	1901	1906	rBV3	117996	170203	4.93%	0.560%
10	14.463	1908	1917	1926	rBV	1415753	2512559	72.78%	8.272%
11	14.993	2004	2007	2012	rVB2	197644	247614	7.17%	0.815%
12	15.257	2049	2052	2056	rVB2	197814	235314	6.82%	0.775%
13	15.322	2061	2063	2067	rVB	172834	190881	5.53%	0.628%
14	16.192	2207	2211	2212	rBV	534967	654665	18.96%	2.155%
15	16.216	2212	2215	2219	rVB	789220	1121304	32.48%	3.692%
16	16.769	2306	2309	2312	rBV	488270	562773	16.30%	1.853%
17	16.987	2343	2346	2349	rBV	621093	743192	21.53%	2.447%
18	17.039	2352	2355	2359	rVB	975110	1347883	39.04%	4.438%
19	17.228	2383	2387	2394	rBV	1547164	2289502	66.32%	7.538%
20	17.298	2396	2399	2407	rVB3	494071	1020950	29.57%	3.361%
21	17.734	2469	2473	2478	rBV	853189	1628723	47.18%	5.362%
22	18.010	2517	2520	2524	rBV2	659381	1034330	29.96%	3.405%
23	18.228	2554	2557	2560	rBV	705233	1019944	29.54%	3.358%
24	18.410	2584	2588	2592	rBV	1336330	2235593	64.76%	7.360%
25	18.851	2660	2663	2667	rBV	1277638	1792102	51.91%	5.900%
26	19.022	2688	2692	2695	rBV	1525988	2179307	63.13%	7.175%
27	19.233	2725	2728	2733	rBV2	1316854	2181444	63.19%	7.182%
28	21.233	3065	3068	3075	rVB	2508316	3452331	100.00%	11.366%

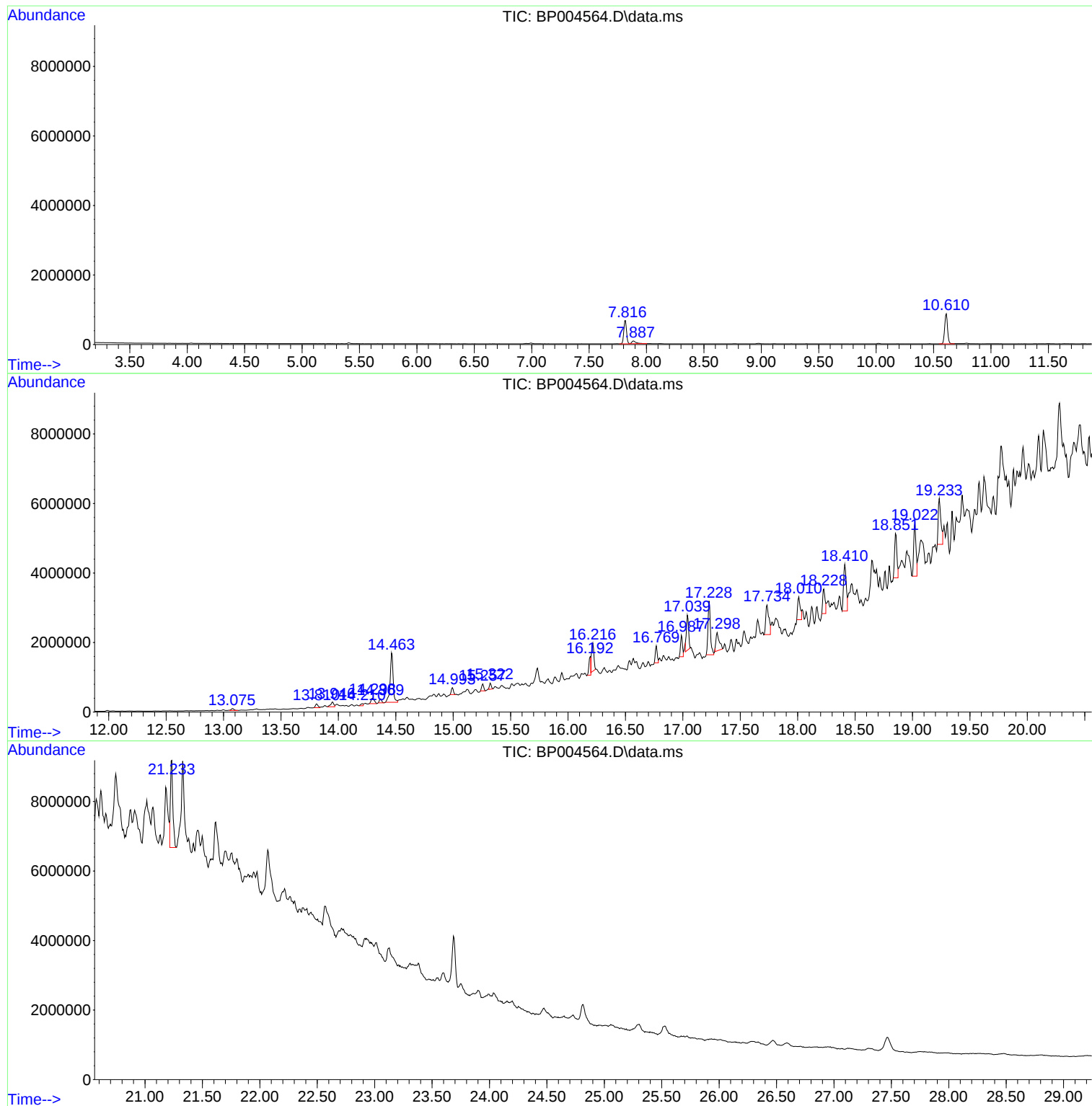
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ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
1717-1718

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP011221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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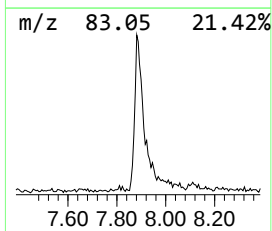
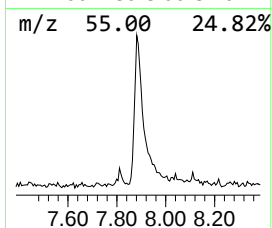
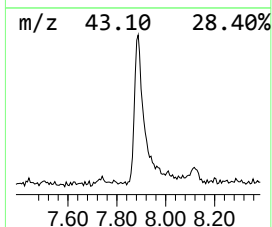
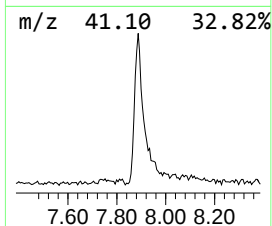
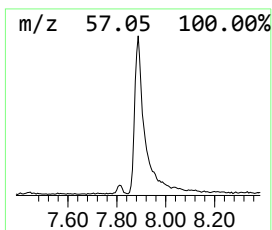
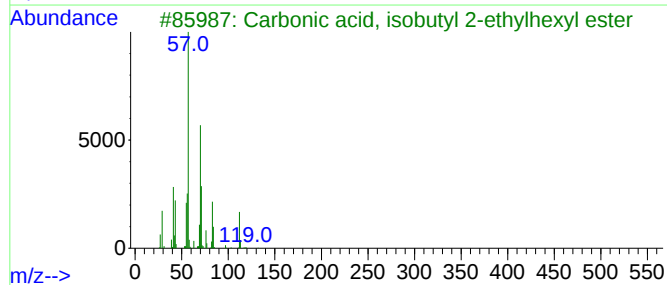
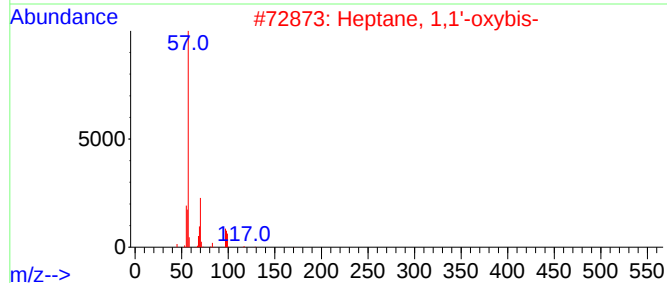
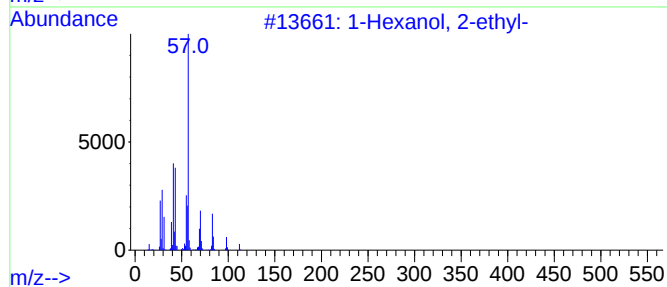
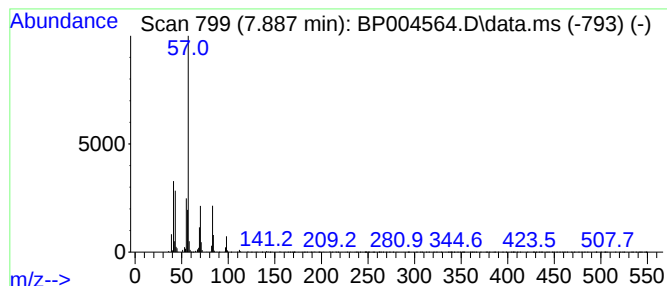
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.887	3.82 ng	210799	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2			Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	59
3			Carbonic acid, isobutyl 2-ethylh...	230	C13H26O3	1000357-82-9	53
4			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	50
5			2-Ethyl-1-hexanol, methyl ether	144	C9H20O	1000365-10-2	50



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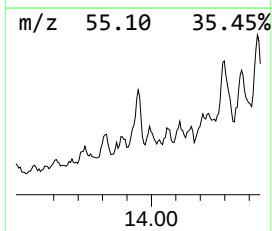
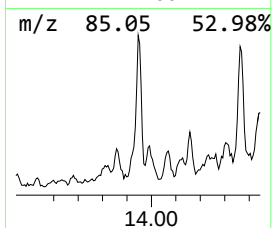
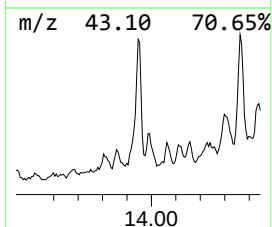
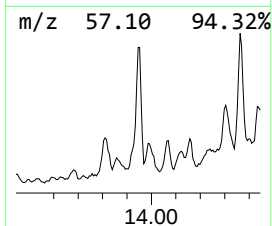
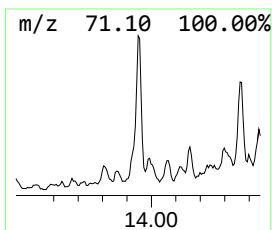
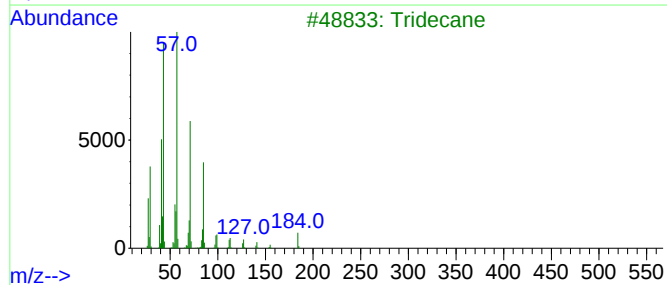
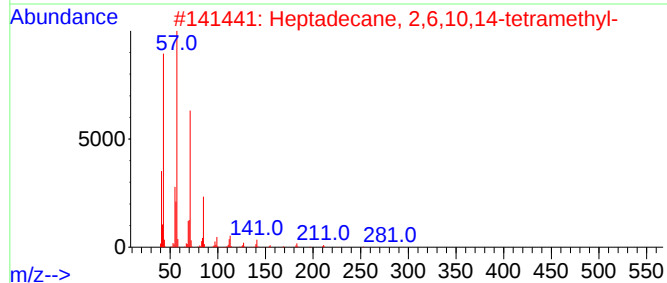
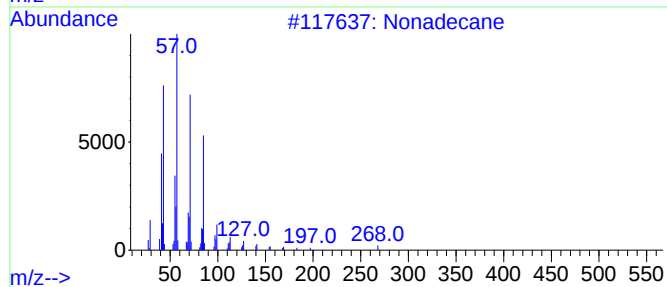
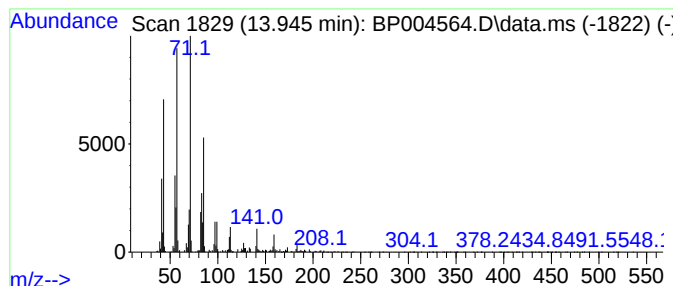
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Nonadecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.945	2.10 ng	263685	Acenaphthene-d10	14.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	86
2			Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	81
3			Tridecane	184	C13H28	000629-50-5	81
4			Decane, 2,3,7-trimethyl-	184	C13H28	062238-13-5	80
5			Hexadecane	226	C16H34	000544-76-3	70



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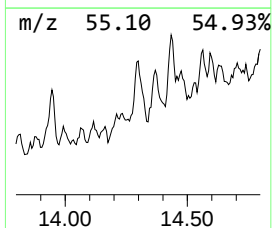
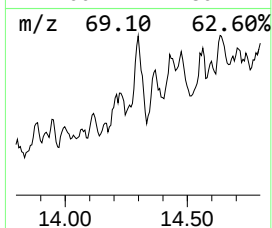
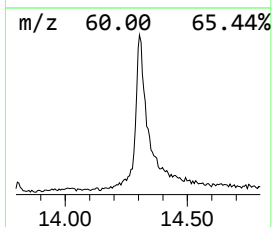
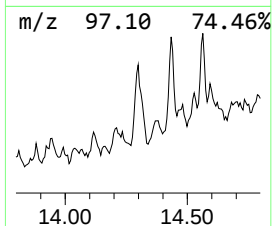
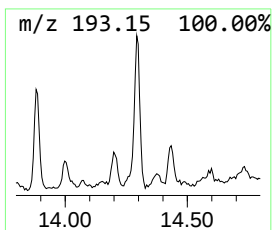
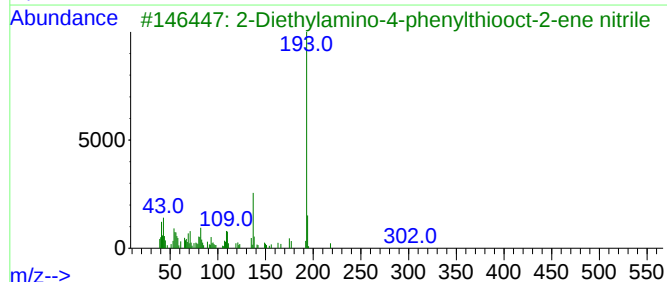
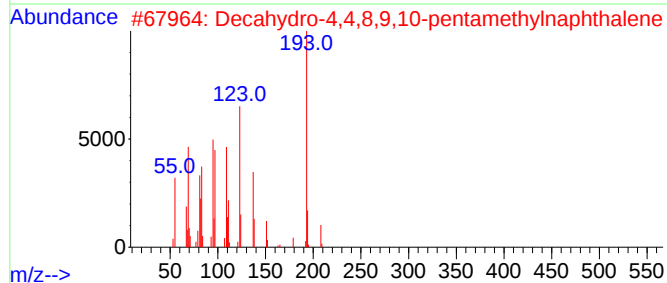
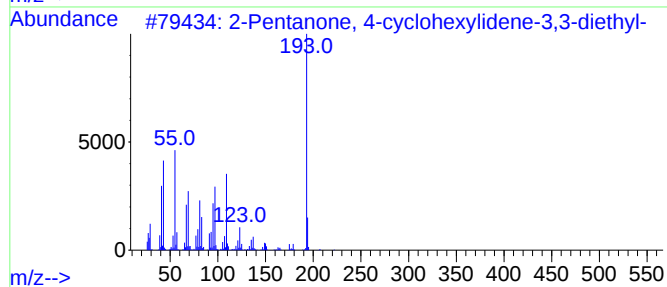
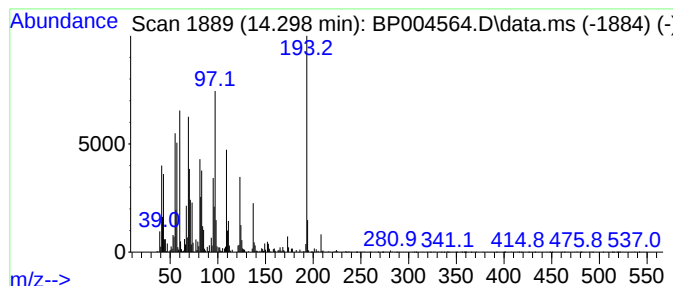
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown14.298 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.298	2.47 ng	310438	Acenaphthene-d10	14.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	43		
2	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	38		
3	2-Diethylamino-4-phenylthiooct-2...	302	C18H26N2S	1000090-57-0	20		
4	Cyclohexane, 1-(cyclohexylmethyl...	194	C14H26	054824-04-3	14		
5	9-Phenanthrenamine	193	C14H11N	000947-73-9	14		



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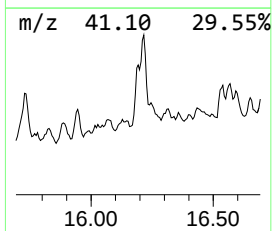
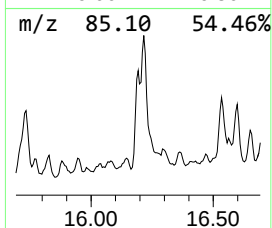
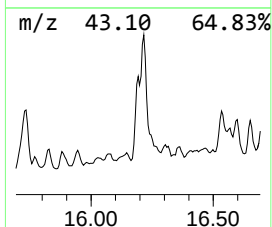
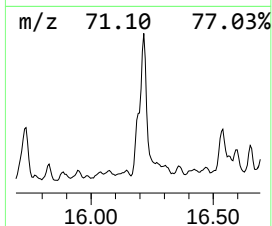
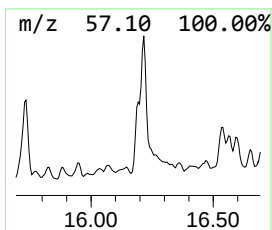
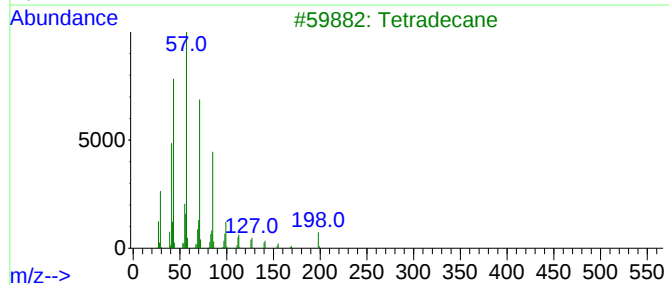
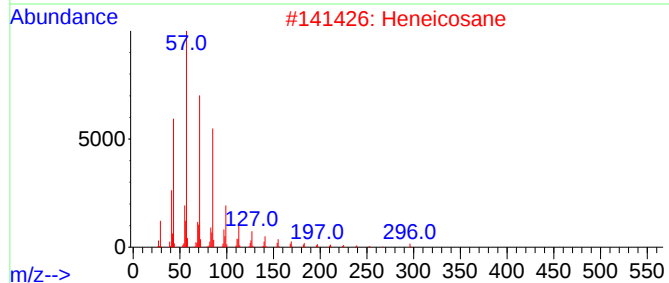
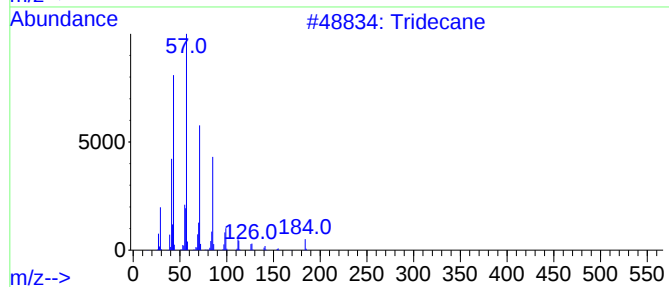
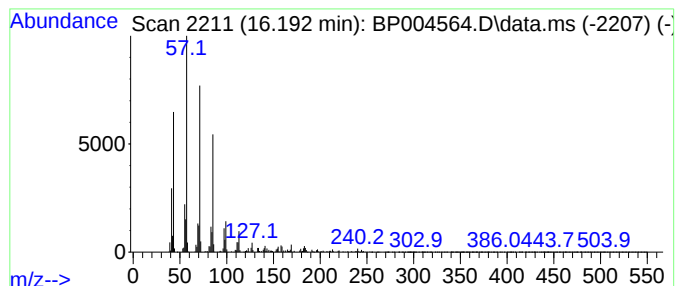
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Peak Number 4 Tridecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.192	5.72 ng	654665	Phenanthrene-d10	17.228

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	93
2			Heneicosane	296	C21H44	000629-94-7	93
3			Tetradecane	198	C14H30	000629-59-4	87
4			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
5			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	83



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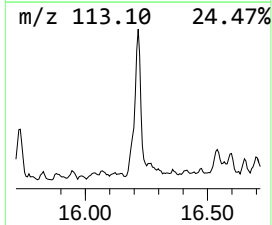
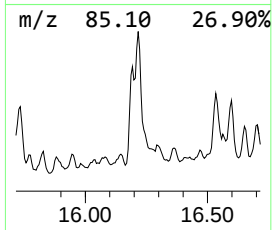
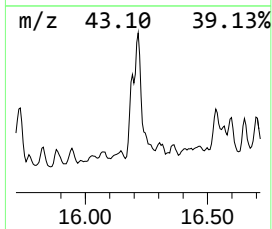
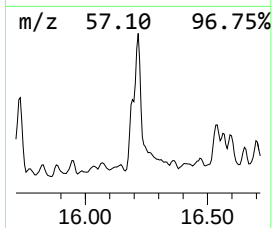
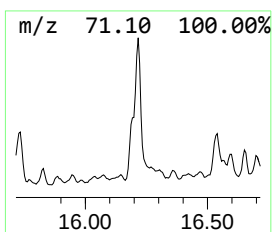
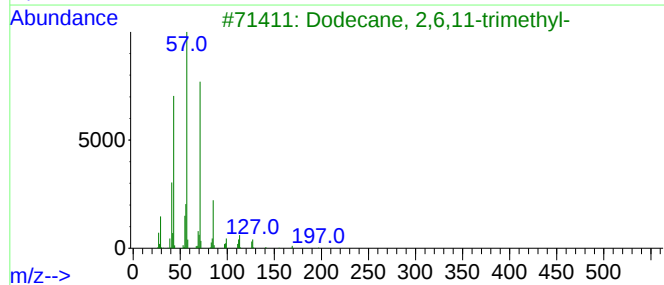
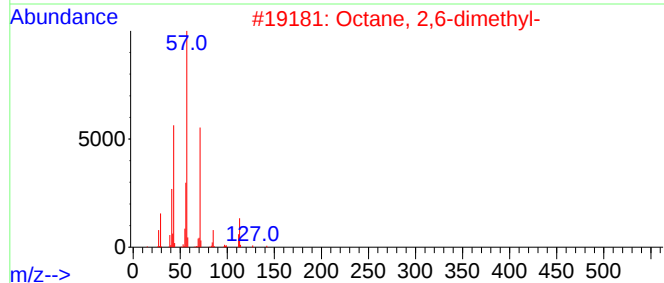
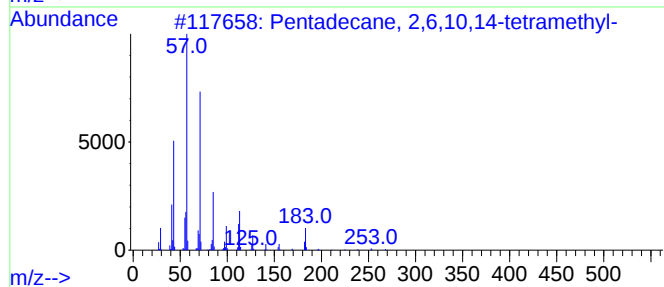
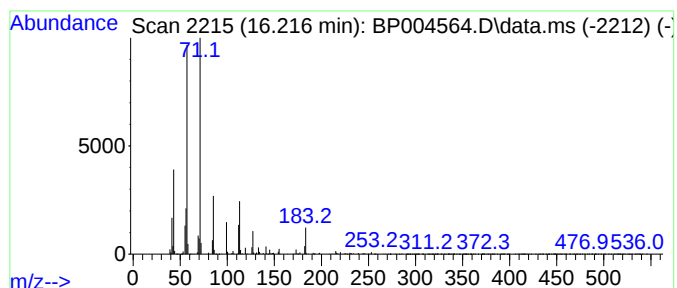
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Pentadecane, 2,6,10,14-tetr... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.216	9.80 ng	1121300	Phenanthrene-d10	17.228

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	91
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
3			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	59
4			Dodecane	170	C12H26	000112-40-3	53
5			Undecane, 3,8-dimethyl-	184	C13H28	017301-30-3	53



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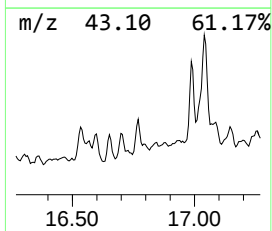
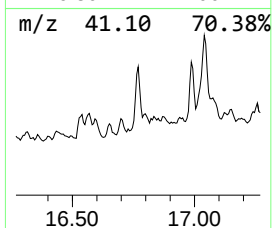
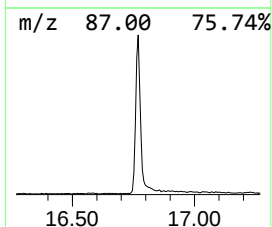
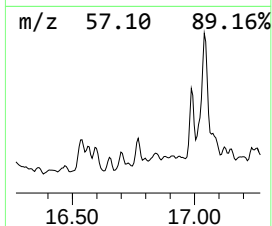
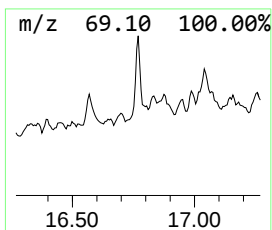
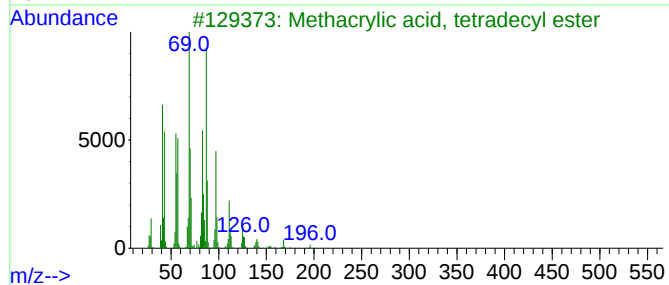
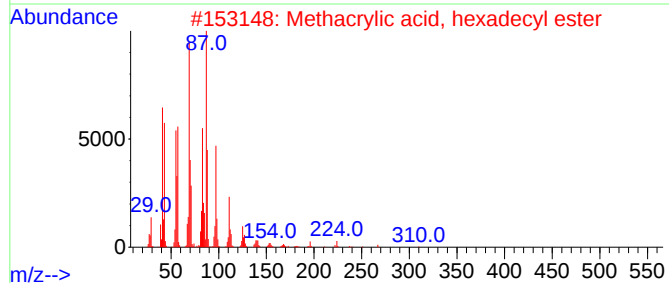
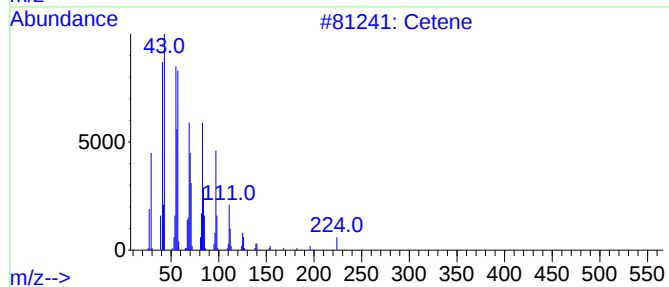
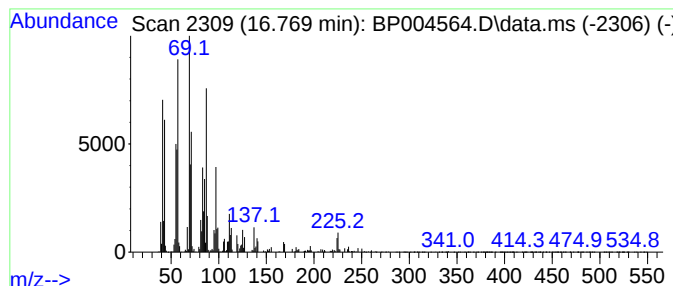
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Cetene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.769	4.92 ng	562773	Phenanthrene-d10	17.228

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cetene	224	C16H32	000629-73-2	80
2			Methacrylic acid, hexadecyl ester	310	C20H38O2	1000340-29-2	80
3			Methacrylic acid, tetradecyl ester	282	C18H34O2	1000340-29-0	74
4			n-Dodecyl methacrylate	254	C16H30O2	000142-90-5	72
5			Methacrylic acid, pentadecyl ester	296	C19H36O2	1000340-29-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011421\
Data File : BP004564.D
Acq On : 14 Jan 2021 19:11
Operator : CG/JU
Sample : M1125-01 10X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
1717-1718

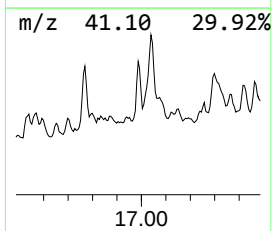
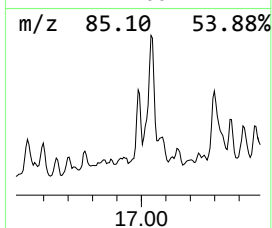
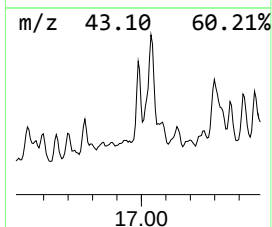
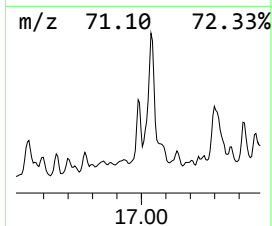
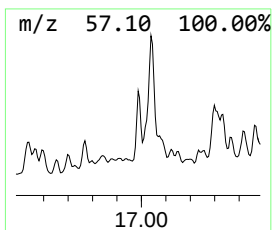
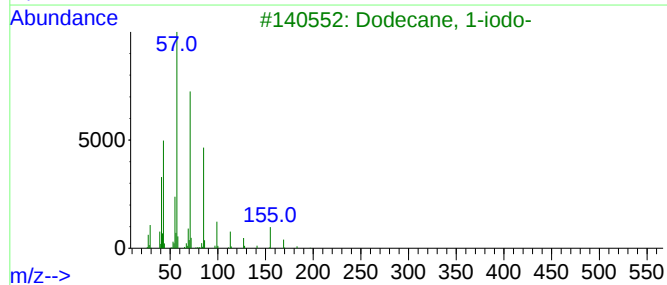
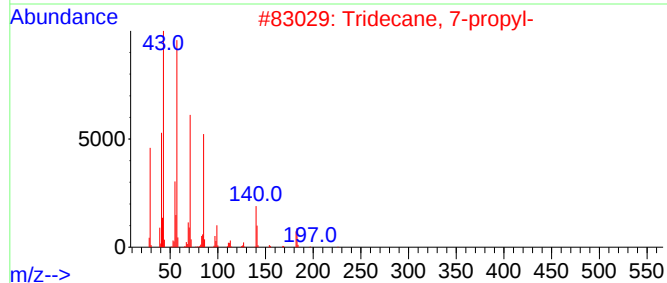
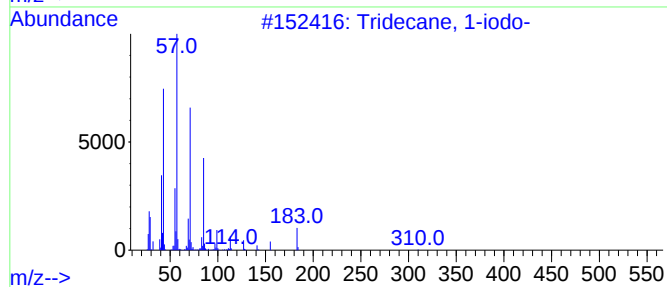
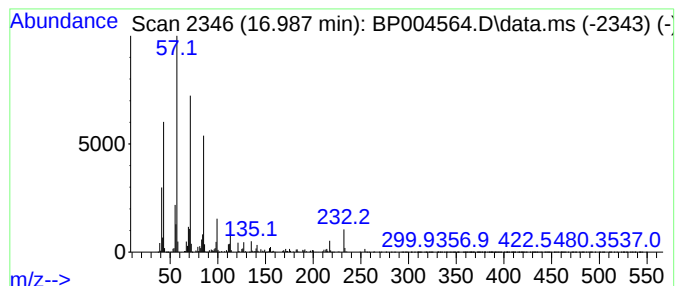
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Tridecane, 1-iodo- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.986	6.49 ng	743192	Phenanthrene-d10	17.228

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	81
2			Tridecane, 7-propyl-	226	C16H34	055045-09-5	81
3			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72
4			Heptadecane	240	C17H36	000629-78-7	72
5			Eicosane	282	C20H42	000112-95-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011421\
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Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
1717-1718

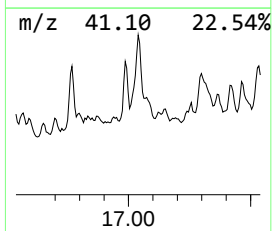
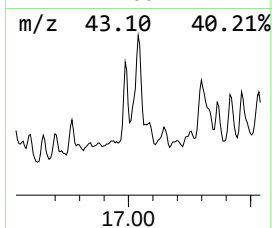
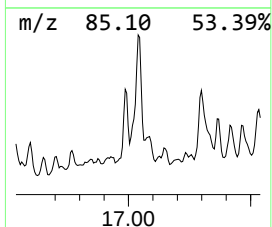
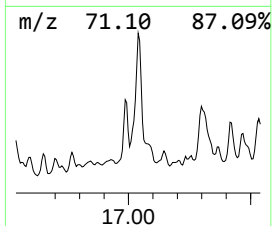
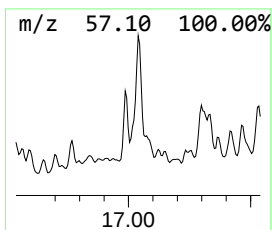
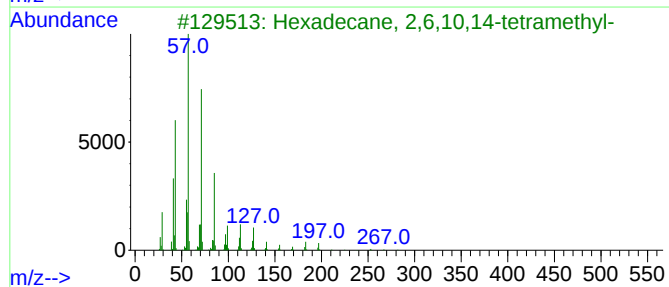
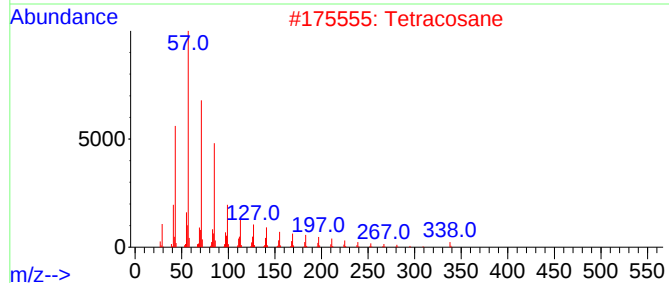
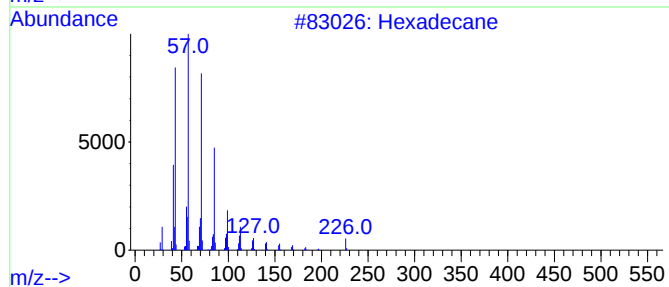
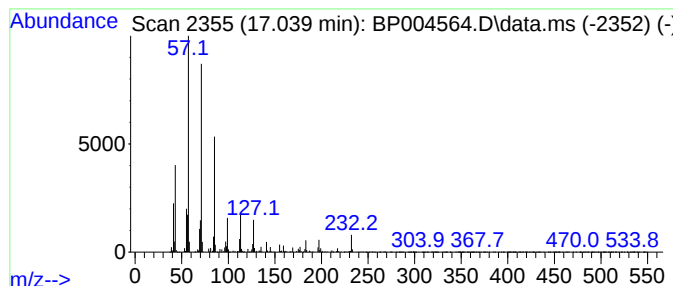
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Hexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.040	11.77 ng	1347880	Phenanthrene-d10	17.228

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	87
2			Tetracosane	338	C24H50	000646-31-1	86
3			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	80
4			Hexacosane	366	C26H54	000630-01-3	78
5			Tetradecane	198	C14H30	000629-59-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011421\
Data File : BP004564.D
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Misc :
ALS Vial : 17 Sample Multiplier: 1

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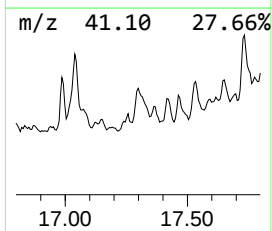
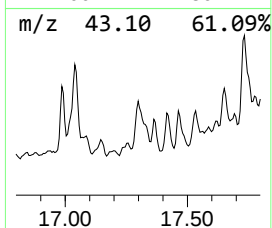
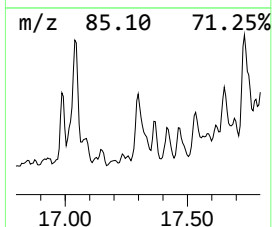
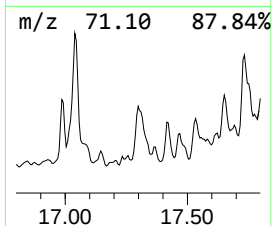
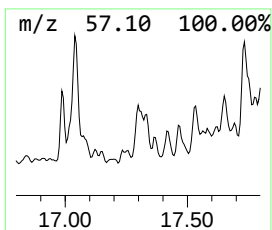
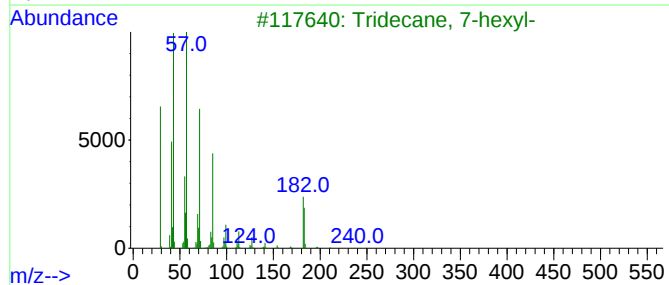
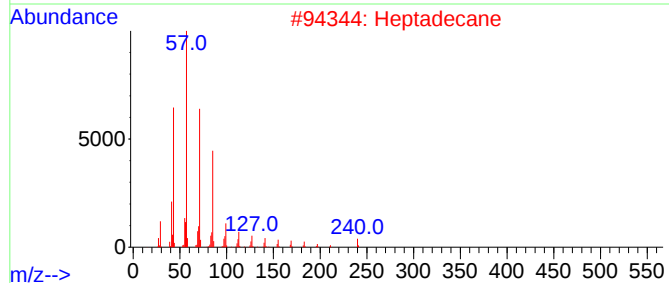
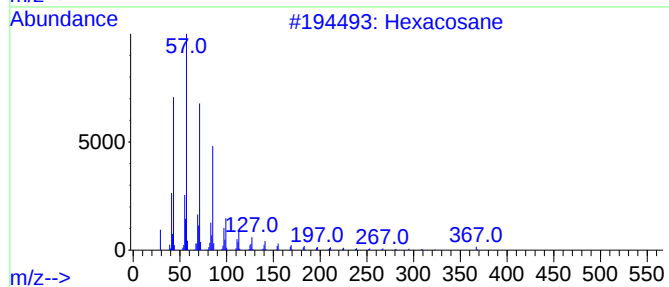
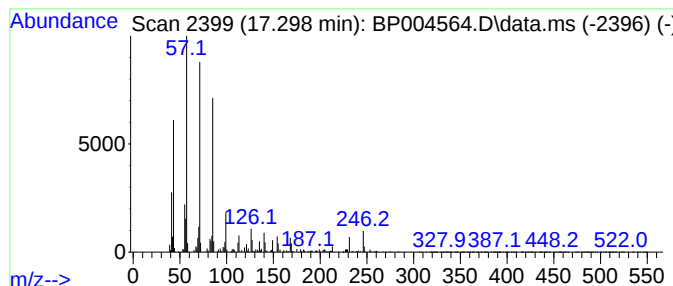
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Hexacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.298	8.92 ng	1020950	Phenanthrene-d10	17.228

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	72
2			Heptadecane	240	C17H36	000629-78-7	72
3			Tridecane, 7-hexyl-	268	C19H40	007225-66-3	72
4			Octacosane	394	C28H58	000630-02-4	64
5			Tridecane, 3-ethyl-	212	C15H32	013286-73-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011421\
Data File : BP004564.D
Acq On : 14 Jan 2021 19:11
Operator : CG/JU
Sample : M1125-01 10X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
1717-1718

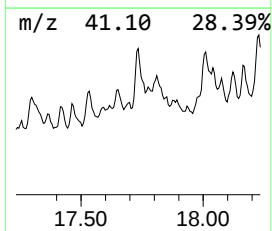
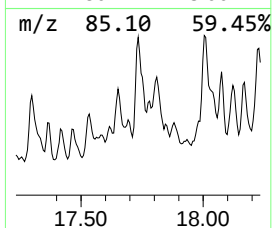
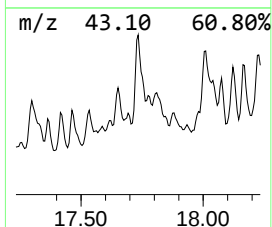
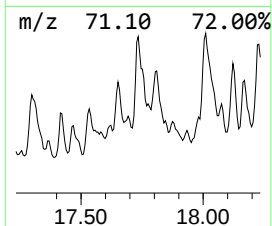
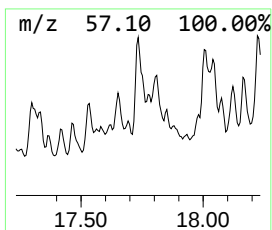
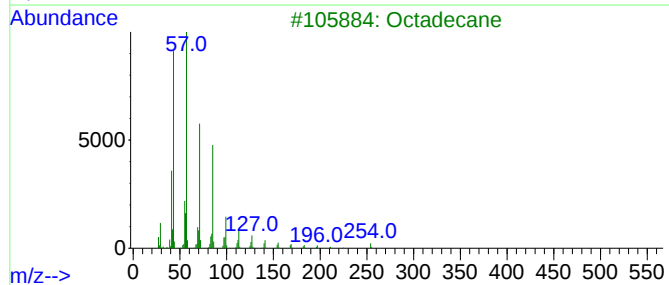
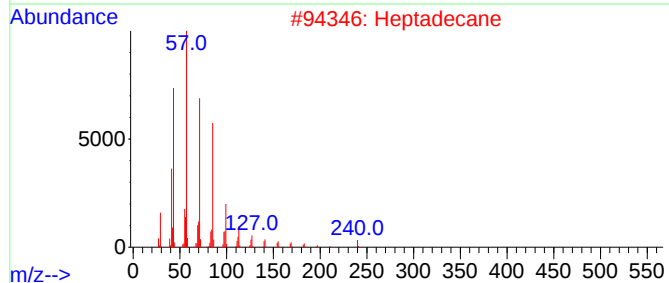
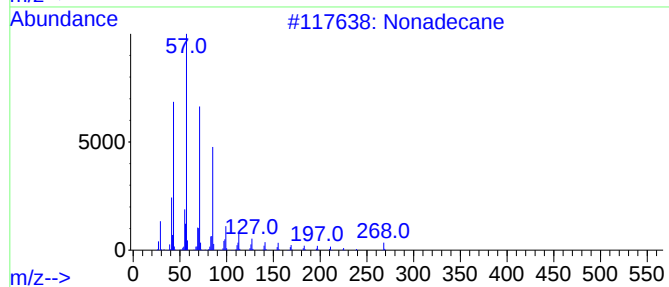
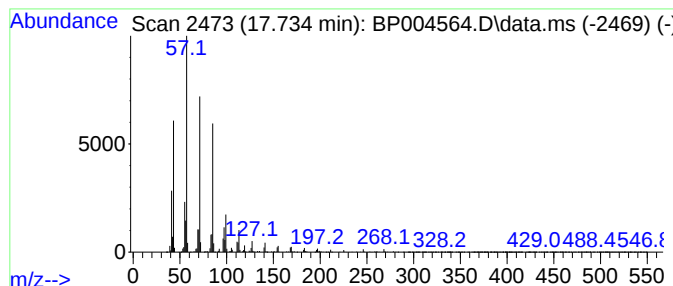
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Heptadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.733	14.23 ng	1628720	Phenanthrene-d10	17.228

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	94
2		Heptadecane	240	C17H36	000629-78-7	91
3		Octadecane	254	C18H38	000593-45-3	87
4		Heneicosane	296	C21H44	000629-94-7	81
5		Octacosane	394	C28H58	000630-02-4	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011421\
Data File : BP004564.D
Acq On : 14 Jan 2021 19:11
Operator : CG/JU
Sample : M1125-01 10X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
1717-1718

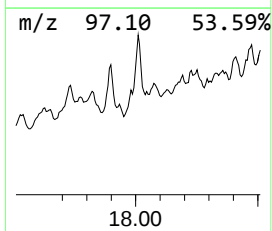
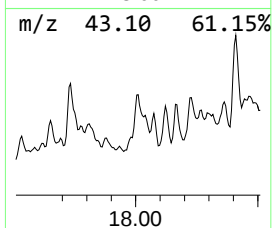
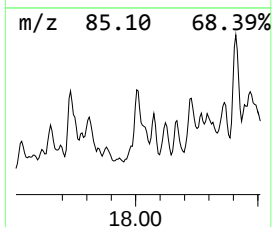
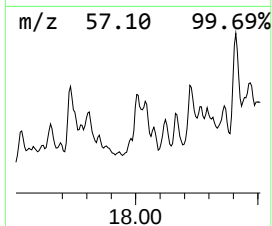
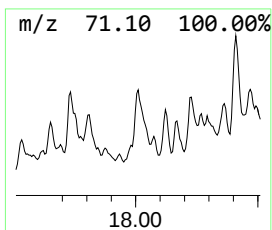
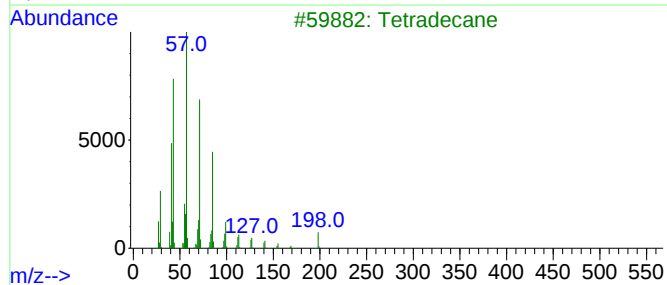
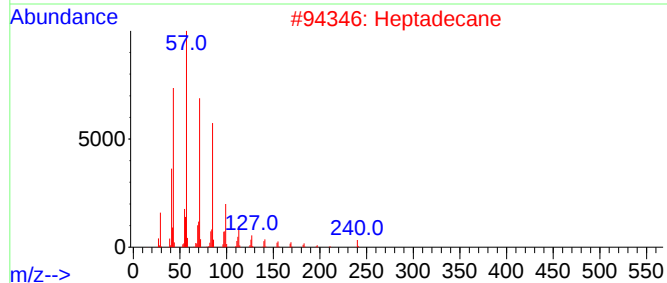
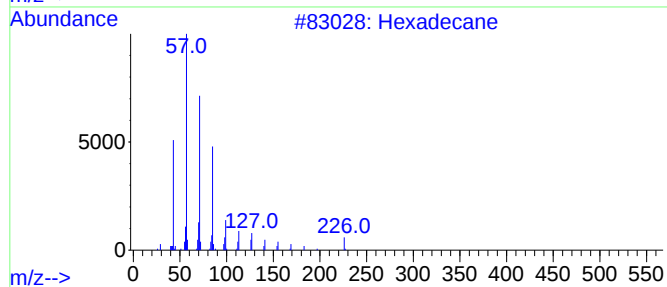
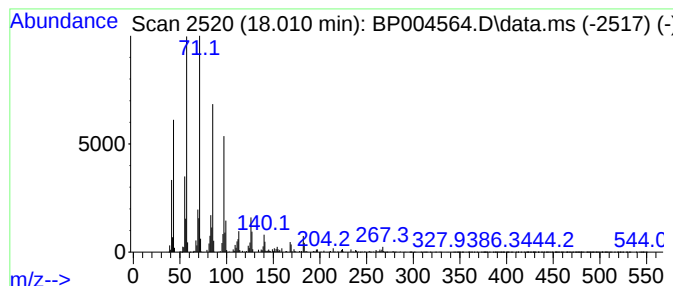
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Tetradecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.010	9.04 ng	1034330	Phenanthrene-d10	17.228

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	74
2			Heptadecane	240	C17H36	000629-78-7	72
3			Tetradecane	198	C14H30	000629-59-4	72
4			Decane, 3-methyl-	156	C11H24	013151-34-3	62
5			Octadecane, 4-methyl-	268	C19H40	010544-95-3	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011421\
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Sample : M1125-01 10X
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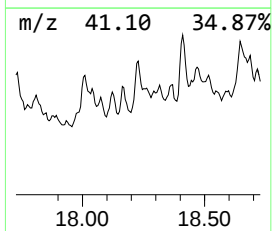
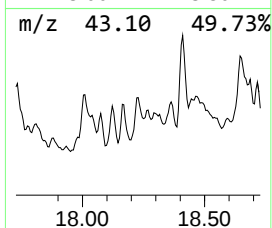
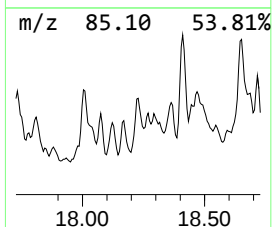
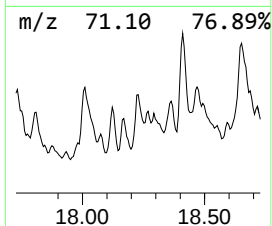
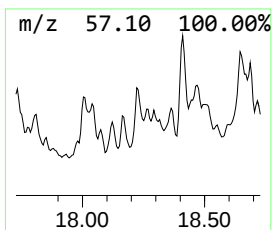
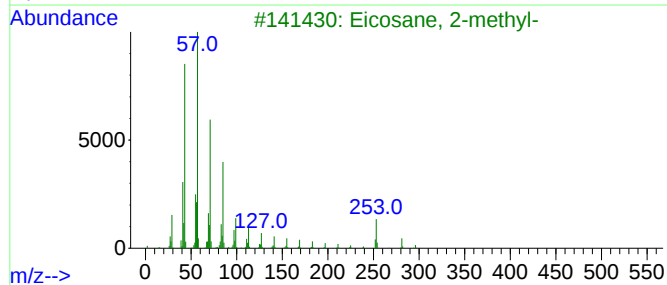
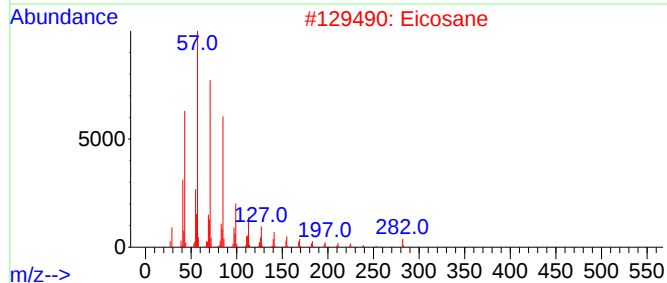
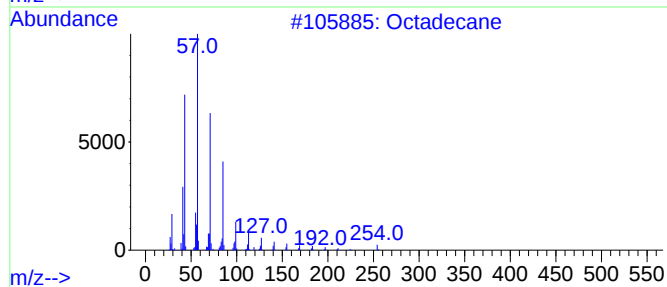
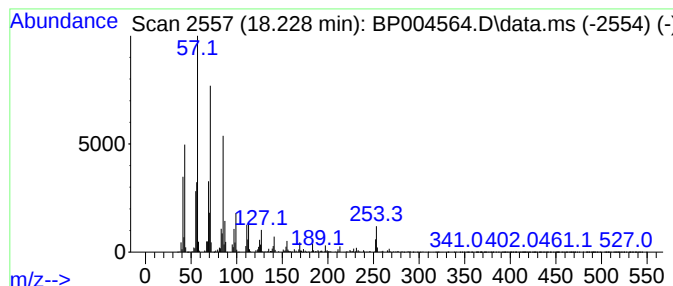
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Octadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.228	8.91 ng	1019940	Phenanthrene-d10		17.228	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octadecane		254	C18H38	000593-45-3	95
2	Eicosane		282	C20H42	000112-95-8	87
3	Eicosane, 2-methyl-		296	C21H44	001560-84-5	86
4	Octacosane		394	C28H58	000630-02-4	83
5	Hentriacontane		437	C31H64	000630-04-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011421\
Data File : BP004564.D
Acq On : 14 Jan 2021 19:11
Operator : CG/JU
Sample : M1125-01 10X
Misc :
ALS Vial : 17 Sample Multiplier: 1

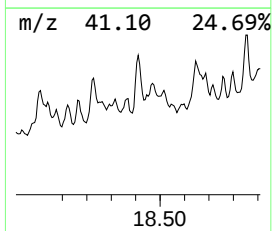
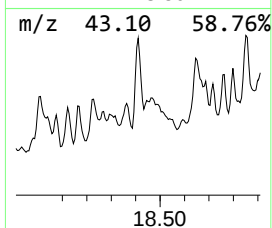
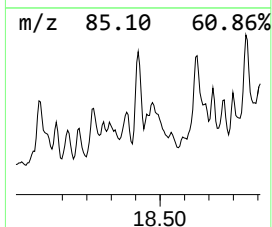
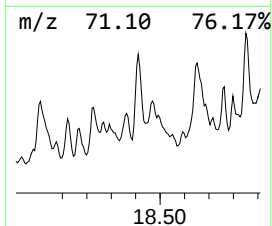
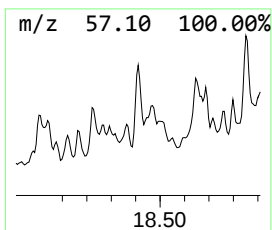
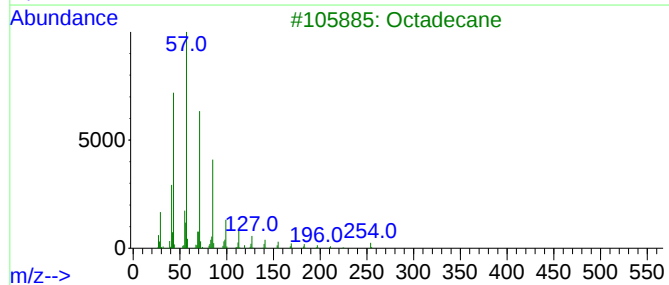
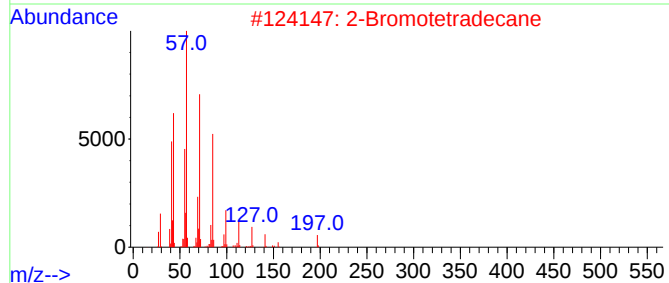
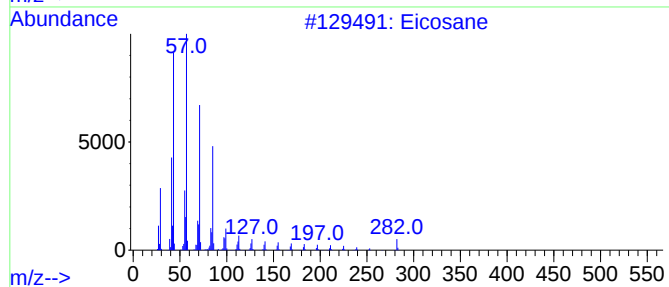
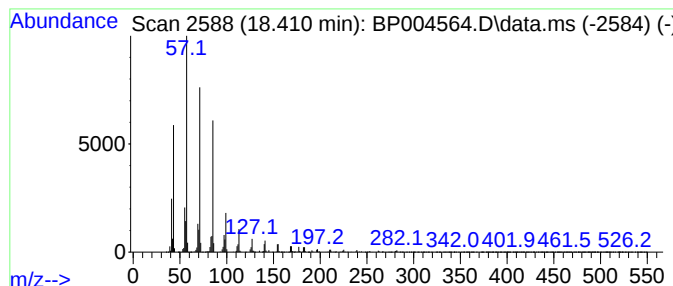
Instrument :
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ClientSampleId :
1717-1718

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP011221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Eicosane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.410	19.53 ng	2235590	Phenanthrene-d10		17.228	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	90
2	2-Bromotetradecane		276	C14H29Br	074036-95-6	90
3	Octadecane		254	C18H38	000593-45-3	90
4	Docosane		310	C22H46	000629-97-0	90
5	Nonadecane, 9-methyl-		282	C20H42	013287-24-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011421\
Data File : BP004564.D
Acq On : 14 Jan 2021 19:11
Operator : CG/JU
Sample : M1125-01 10X
Misc :
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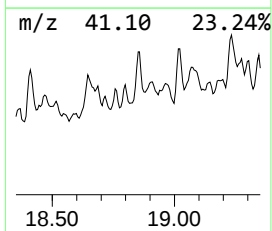
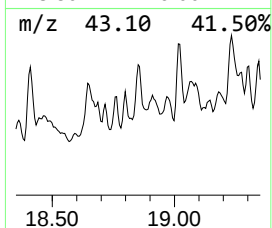
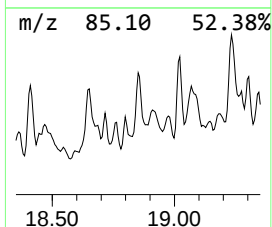
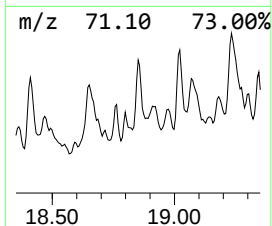
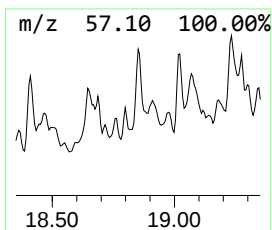
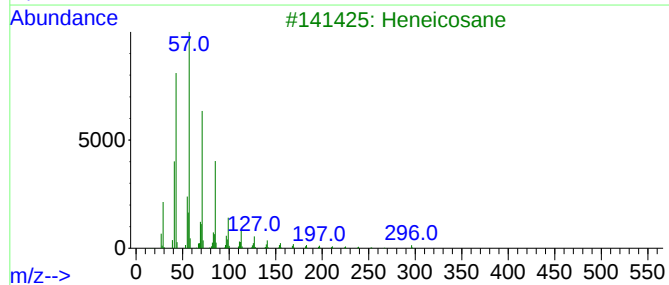
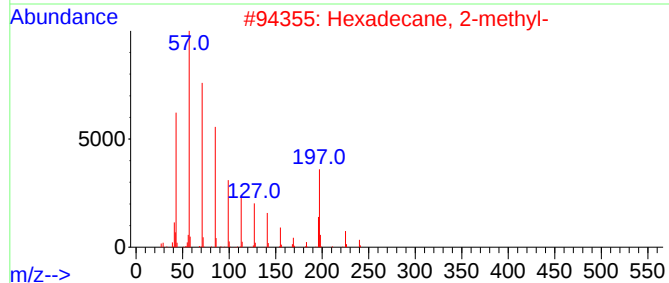
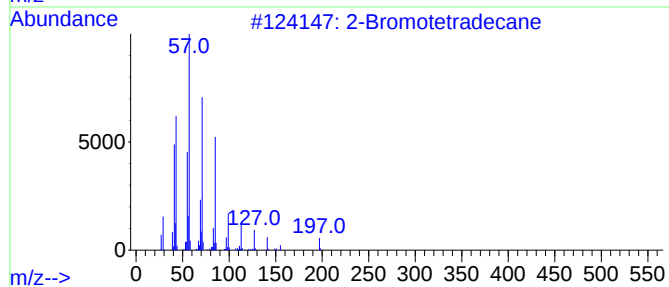
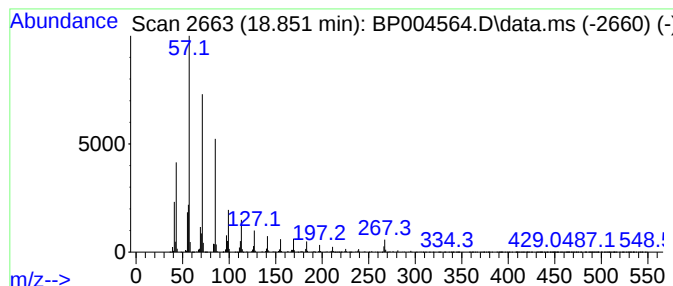
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ClientSampleId :
1717-1718

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP011221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 2-Bromotetradecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.851	15.65 ng	1792100	Phenanthrene-d10		17.228	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Bromotetradecane		276	C14H29Br	074036-95-6	91
2	Hexadecane, 2-methyl-		240	C17H36	001560-92-5	90
3	Heneicosane		296	C21H44	000629-94-7	90
4	Decane, 3,8-dimethyl-		170	C12H26	017312-55-9	81
5	Eicosane, 10-methyl-		296	C21H44	054833-23-7	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011421\
Data File : BP004564.D
Acq On : 14 Jan 2021 19:11
Operator : CG/JU
Sample : M1125-01 10X
Misc :
ALS Vial : 17 Sample Multiplier: 1

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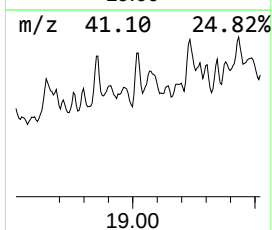
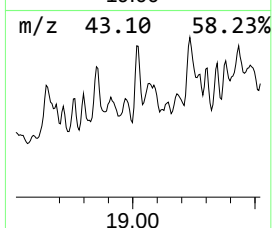
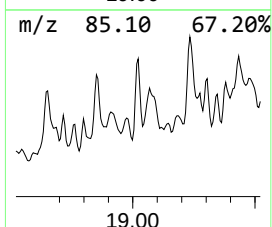
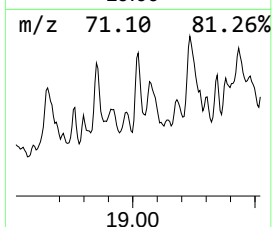
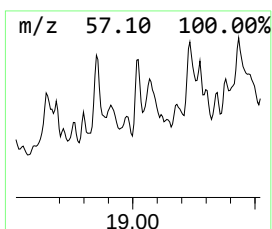
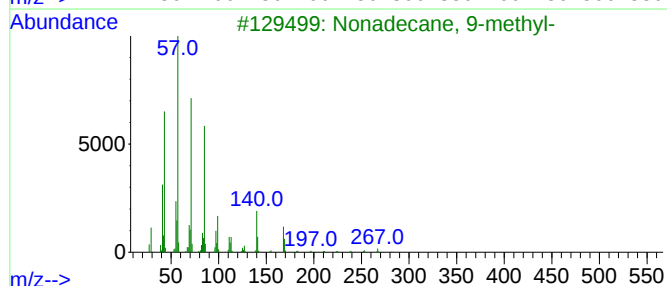
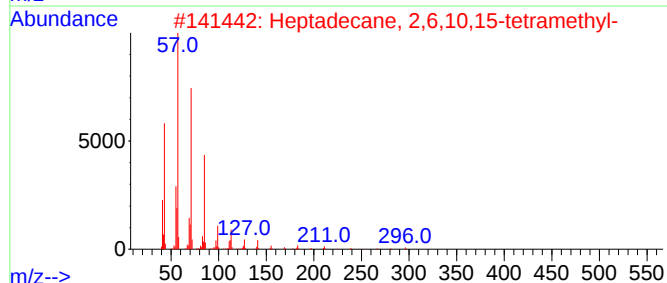
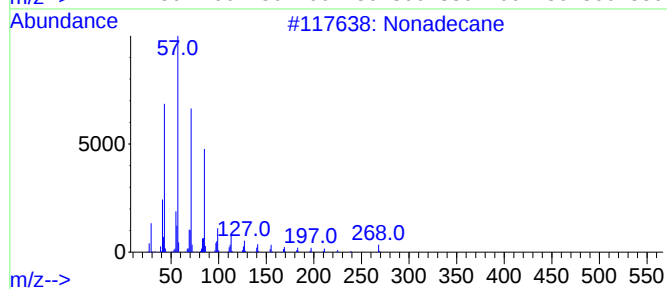
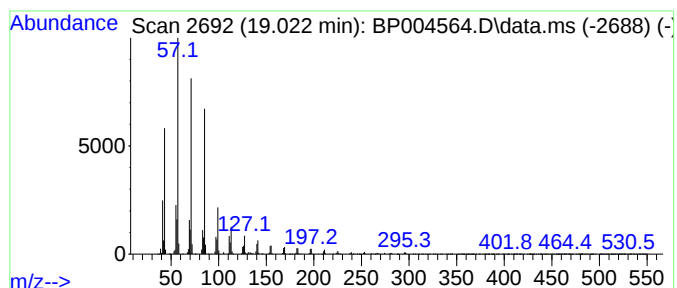
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Heptadecane, 2,6,10,15-tetr... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.022	19.04 ng	2179310	Phenanthrene-d10	17.228

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	95
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
3		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	93
4		Heneicosane	296	C21H44	000629-94-7	91
5		Eicosane	282	C20H42	000112-95-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011421\
Data File : BP004564.D
Acq On : 14 Jan 2021 19:11
Operator : CG/JU
Sample : M1125-01 10X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
1717-1718

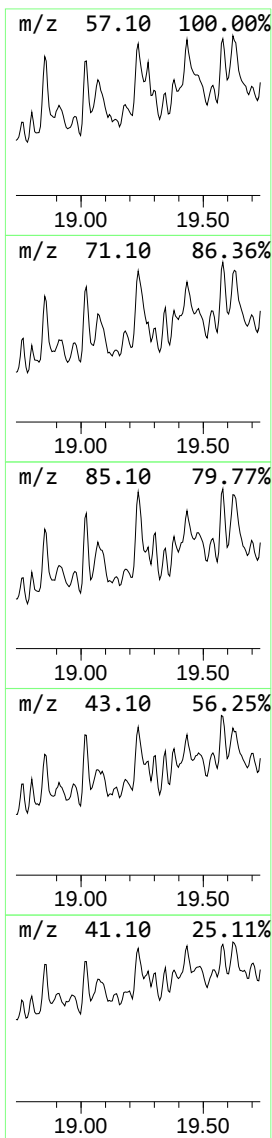
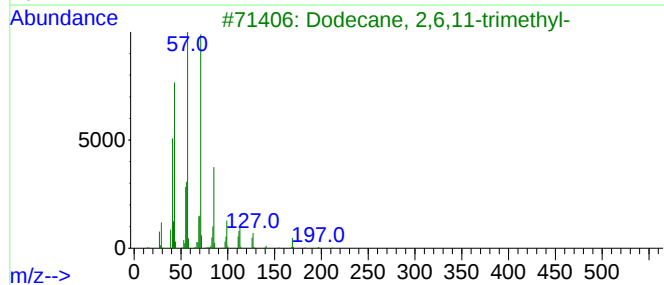
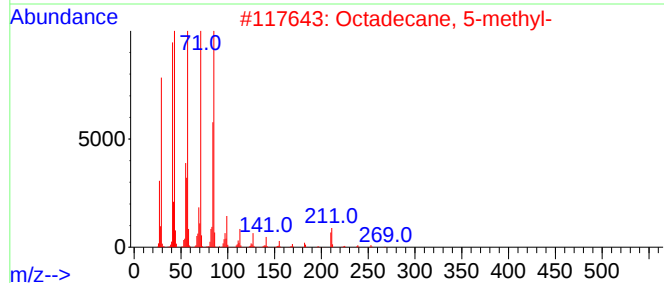
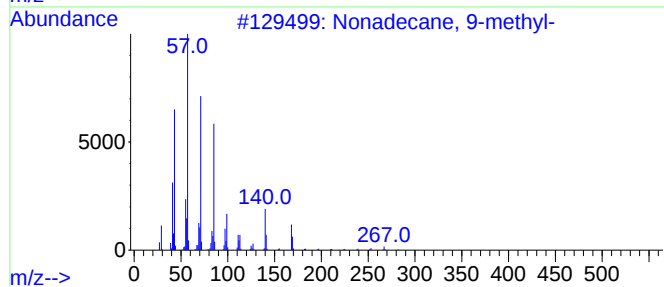
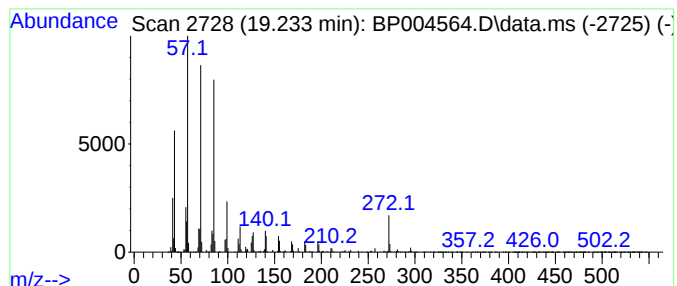
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Nonadecane, 9-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.233	19.06 ng	2181440	Phenanthrene-d10	17.228

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	90
2			Octadecane, 5-methyl-	268	C19H40	025117-35-5	81
3			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	78
4			Heneicosane	296	C21H44	000629-94-7	70
5			3,5-Dimethyldodecane	198	C14H30	107770-99-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011421\
 Data File : BP004564.D
 Acq On : 14 Jan 2021 19:11
 Operator : CG/JU
 Sample : M1125-01 10X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 1717-1718

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP011221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Hexanol, 2-et...	7.887	3.8	ng	210799	1	7.816	1103890	20.0
Nonadecane	13.945	2.1	ng	263685	3	14.463	2512560	20.0
unknown14.298	14.298	2.5	ng	310438	3	14.463	2512560	20.0
Tridecane	16.192	5.7	ng	654665	4	17.228	2289500	20.0
Pentadecane, 2,...	16.216	9.8	ng	1121300	4	17.228	2289500	20.0
Cetene	16.769	4.9	ng	562773	4	17.228	2289500	20.0
Tridecane, 1-iodo-	16.986	6.5	ng	743192	4	17.228	2289500	20.0
Hexadecane	17.040	11.8	ng	1347880	4	17.228	2289500	20.0
Hexacosane	17.298	8.9	ng	1020950	4	17.228	2289500	20.0
Heptadecane	17.733	14.2	ng	1628720	4	17.228	2289500	20.0
Tetradecane	18.010	9.0	ng	1034330	4	17.228	2289500	20.0
Octadecane	18.228	8.9	ng	1019940	4	17.228	2289500	20.0
Eicosane	18.410	19.5	ng	2235590	4	17.228	2289500	20.0
2-Bromotetradecane	18.851	15.7	ng	1792100	4	17.228	2289500	20.0
Heptadecane, 2,...	19.022	19.0	ng	2179310	4	17.228	2289500	20.0
Nonadecane, 9-m...	19.233	19.1	ng	2181440	4	17.228	2289500	20.0