

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072919\
 Data File : BF115834.D
 Acq On : 29 Jul 2019 16:14
 Operator : HP/JU
 Sample : K4036-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WS-02

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF071219.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.134	3	8	37	rVB	2245969	3356113	73.86%	7.960%
2	5.075	504	508	526	rVB	49372	75415	1.66%	0.179%
3	5.481	573	577	600	rBV	2097124	2394263	52.69%	5.679%
4	6.487	744	748	768	rBV	1973017	1954184	43.01%	4.635%
5	6.634	768	773	776	rBV	3347708	3134627	68.99%	7.435%
6	6.857	808	811	815	rBV	949310	819387	18.03%	1.943%
7	7.016	834	838	842	rBV	3012746	2741200	60.33%	6.502%
8	7.422	902	907	910	rBV	2290204	2200017	48.42%	5.218%
9	8.139	1025	1029	1047	rBV	1162127	1042034	22.93%	2.471%
10	9.216	1207	1212	1215	rBV	4326754	4053114	89.20%	9.613%
11	9.604	1275	1278	1283	rBV	116214	98084	2.16%	0.233%
12	9.892	1323	1327	1343	rVB	1471867	1257784	27.68%	2.983%
13	10.680	1457	1461	1478	rBV	2243546	2119623	46.65%	5.027%
14	11.374	1576	1579	1588	rBV	1304768	1292600	28.45%	3.066%
15	11.898	1663	1668	1683	rBV2	252186	395839	8.71%	0.939%
16	12.592	1782	1786	1793	rBV4	22486	54473	1.20%	0.129%
17	12.686	1796	1802	1815	rVV	426472	702571	15.46%	1.666%
18	12.963	1844	1849	1852	rBV	4449451	4543658	100.00%	10.777%
19	13.857	1998	2001	2003	rBV	88625	78419	1.73%	0.186%
20	13.898	2003	2008	2021	rVV4	75207	195509	4.30%	0.464%
21	14.015	2024	2028	2037	rVB	1320404	1216855	26.78%	2.886%
22	14.174	2051	2055	2063	rBV	235870	218707	4.81%	0.519%
23	14.480	2102	2107	2114	rBV	498159	472292	10.39%	1.120%
24	14.786	2154	2159	2166	rBV	841440	859694	18.92%	2.039%
25	15.121	2211	2216	2225	rBV	1052253	1259258	27.71%	2.987%
26	15.504	2272	2281	2295	rBV2	1247667	2668472	58.73%	6.329%
27	15.762	2323	2325	2331	rVB	39289	53278	1.17%	0.126%
28	15.933	2347	2354	2367	rBV	955713	1581488	34.81%	3.751%
29	16.439	2432	2440	2456	rBV	505698	972062	21.39%	2.306%
30	17.027	2533	2540	2548	rVB	139285	283061	6.23%	0.671%
31	17.721	2655	2658	2667	rVB3	33002	67996	1.50%	0.161%

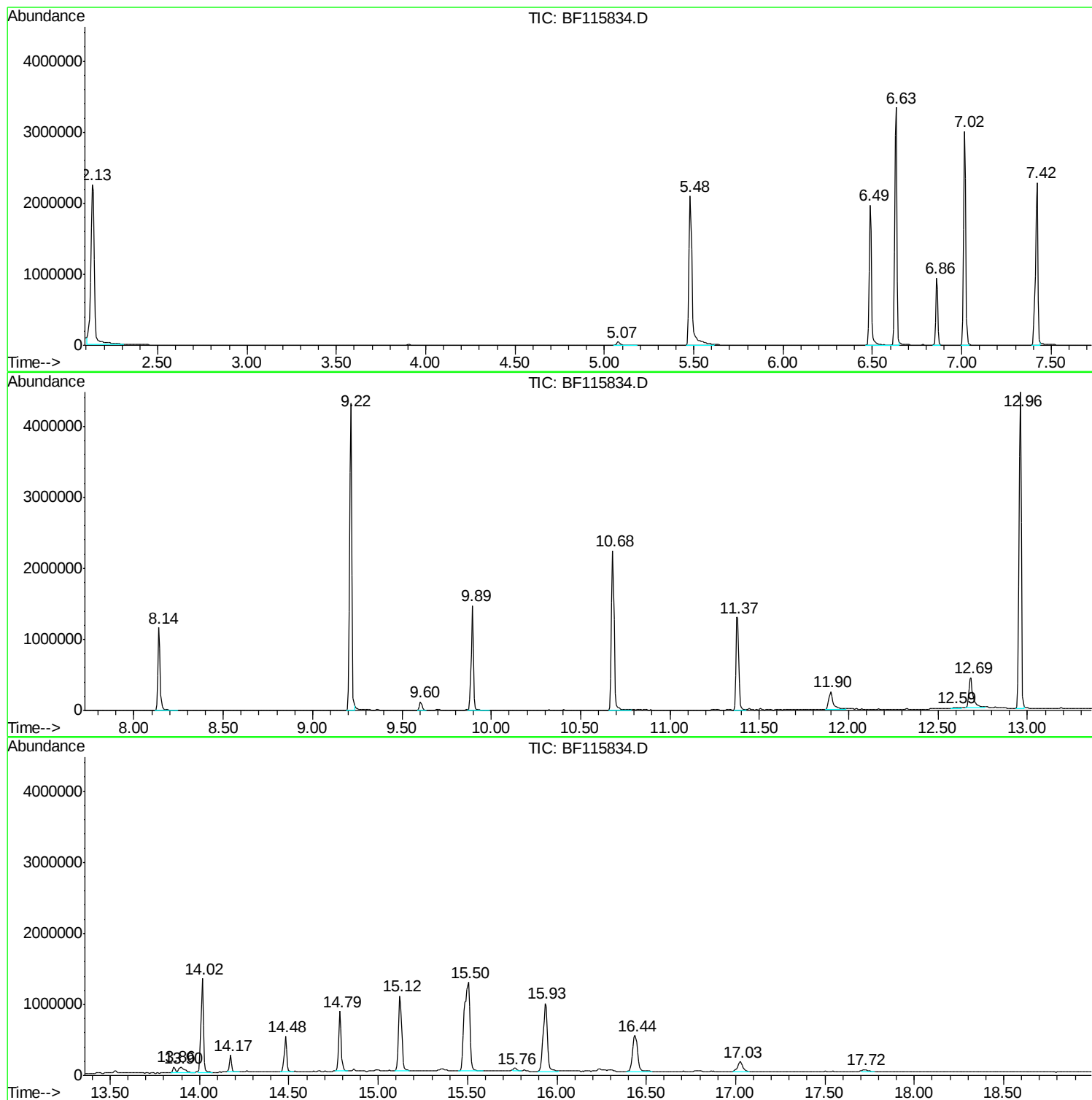
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ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WS-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.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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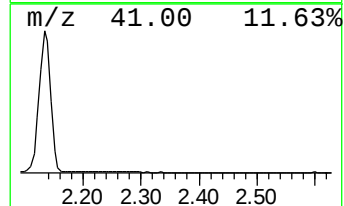
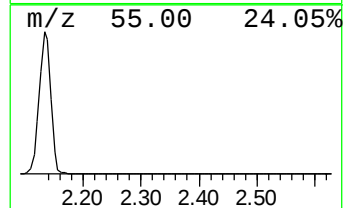
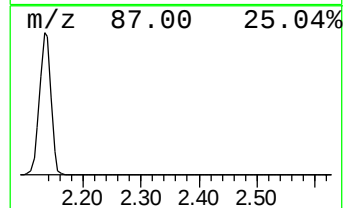
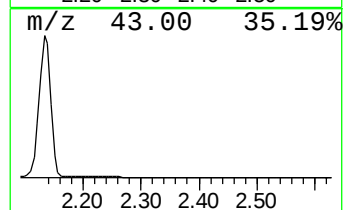
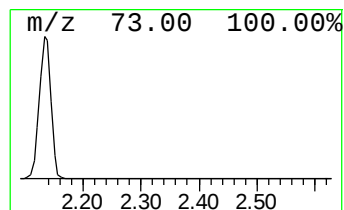
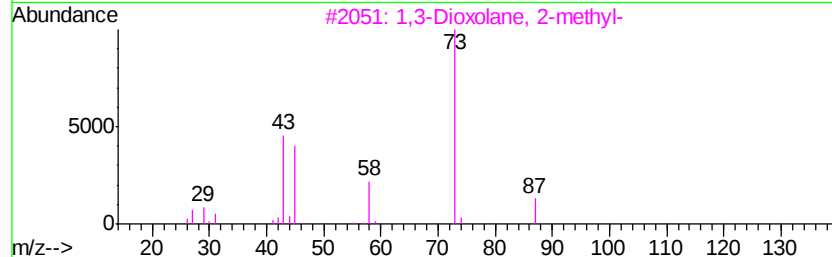
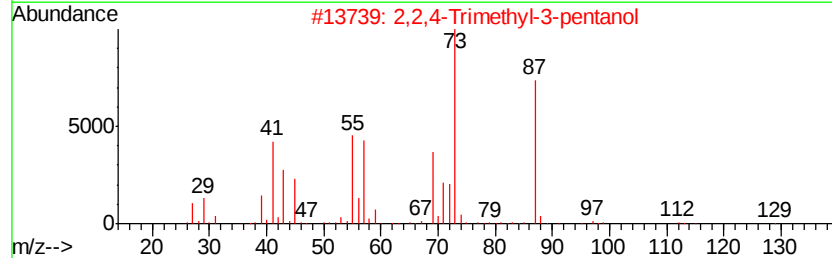
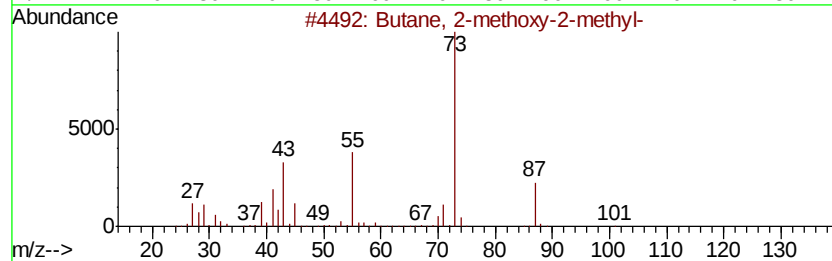
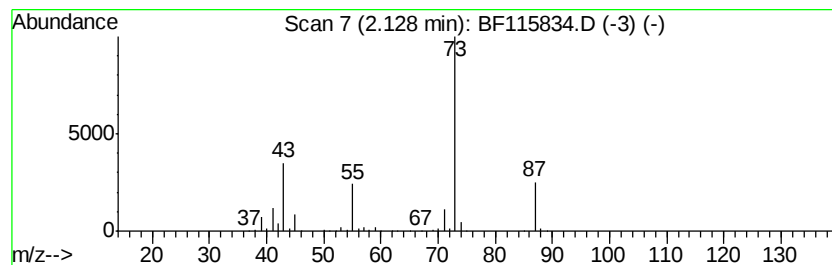
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.13	81.92 ng	3356110	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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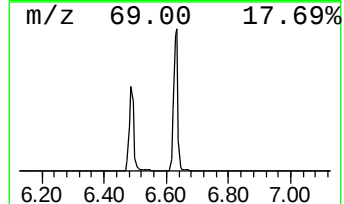
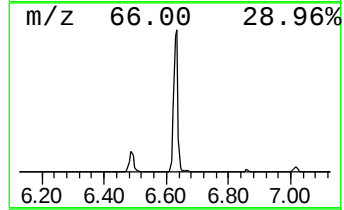
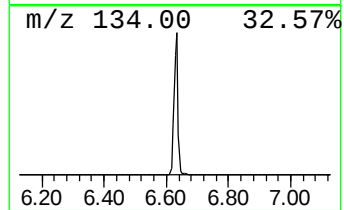
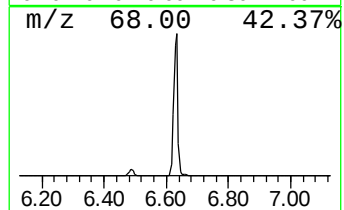
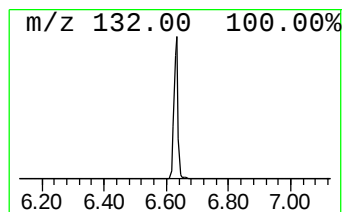
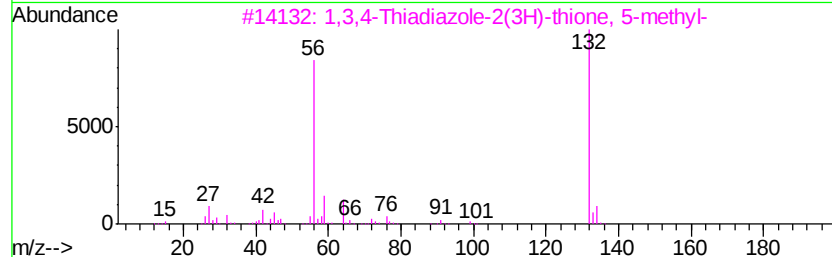
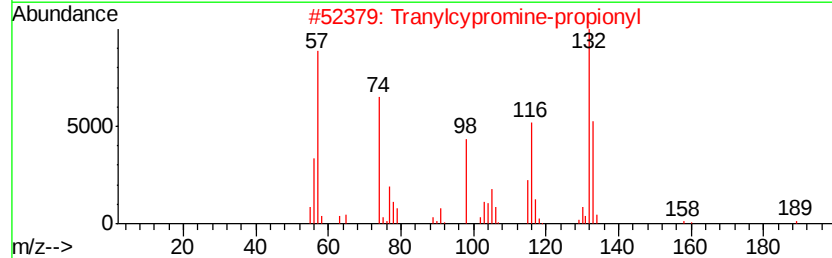
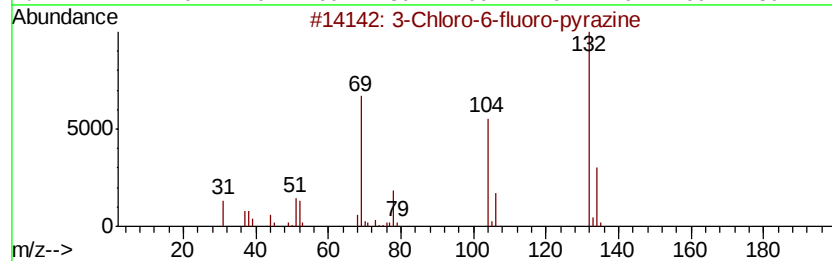
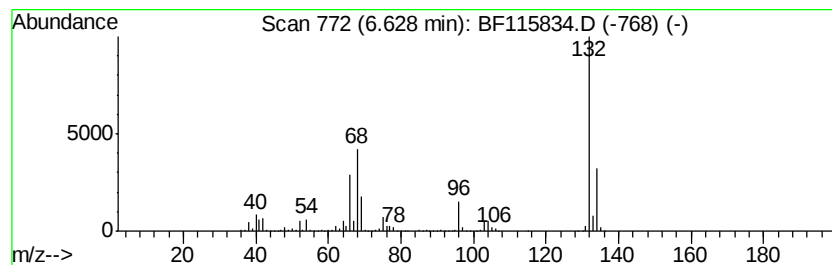
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown6.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	76.51 ng	3134630	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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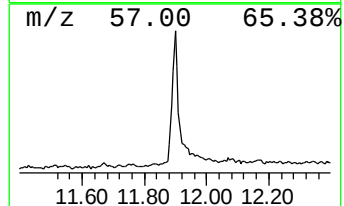
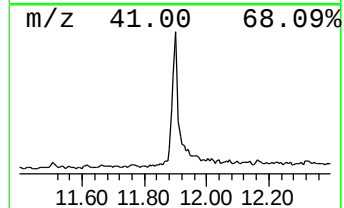
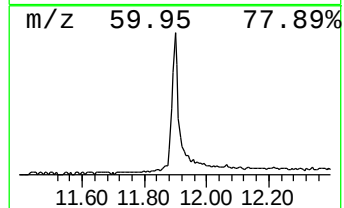
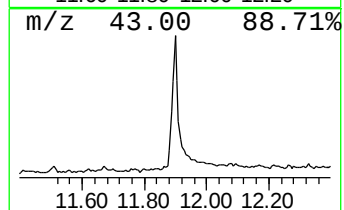
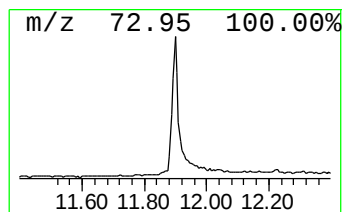
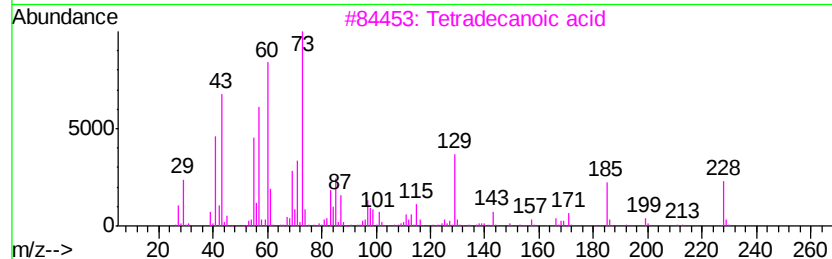
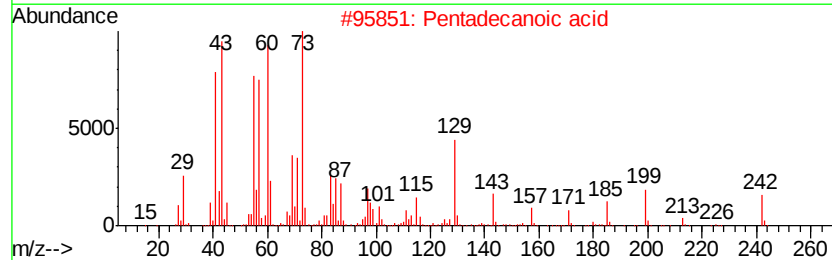
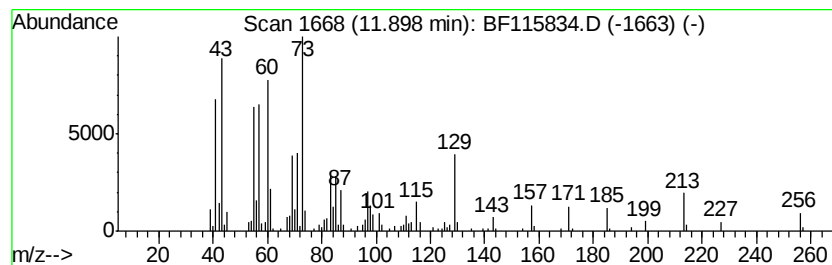
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	6.12 ng	395839	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4		Tridecanoic acid	214	C13H26O2	000638-53-9	83
5		n-Decanoic acid	172	C10H20O2	000334-48-5	76



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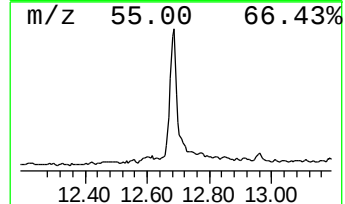
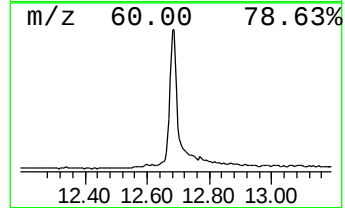
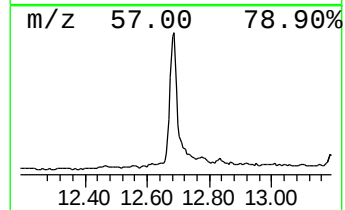
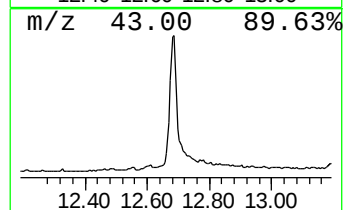
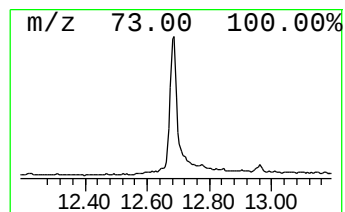
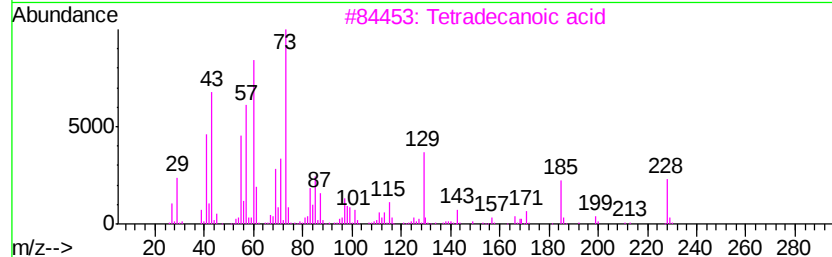
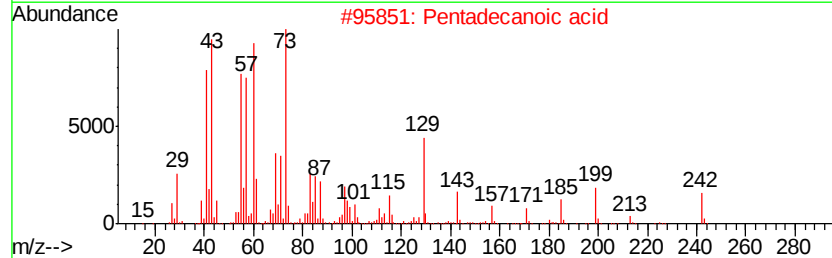
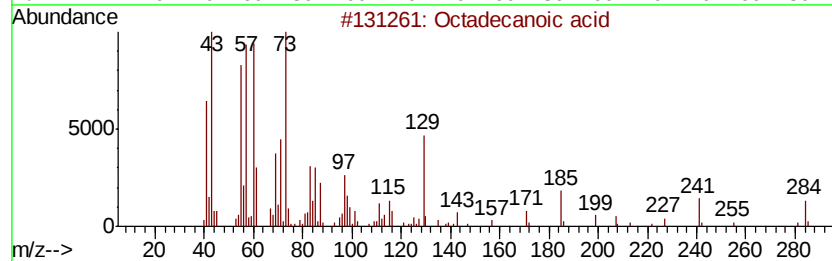
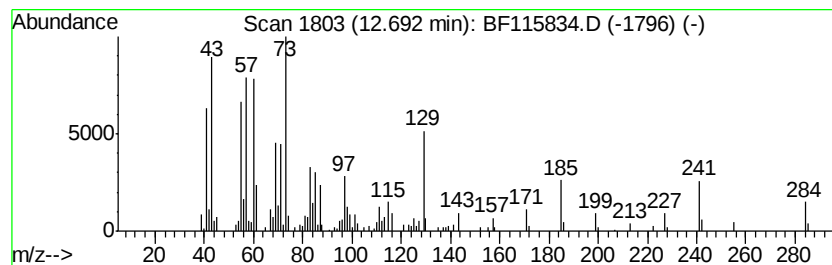
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Peak Number 5 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.69	10.87 ng	702571	Phenanthrene-d10	11.37

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	87
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	81
5		Undecanoic acid	186	C11H22O2	000112-37-8	81



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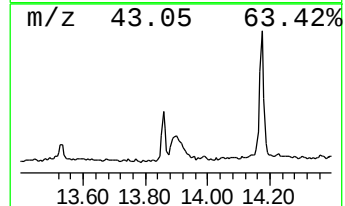
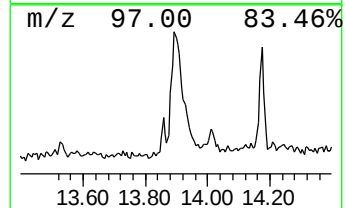
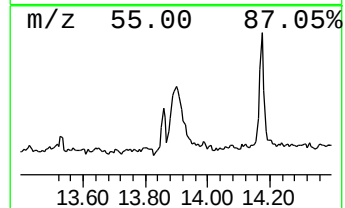
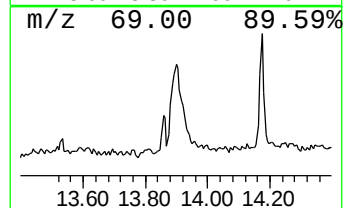
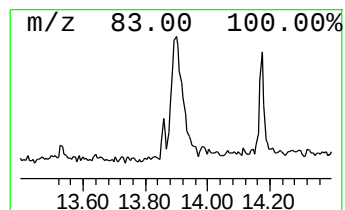
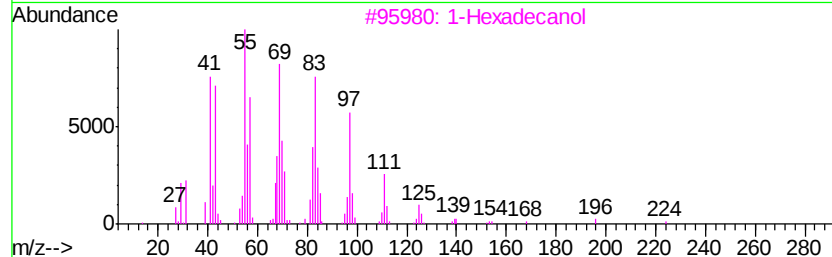
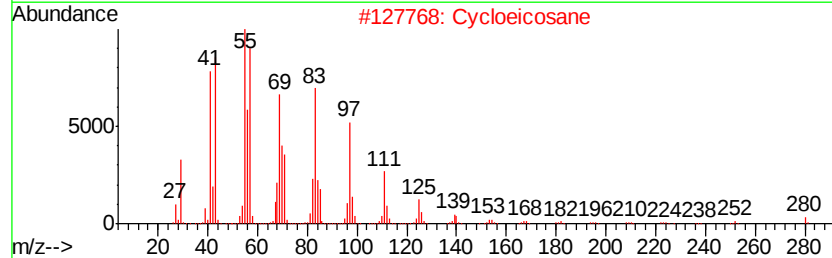
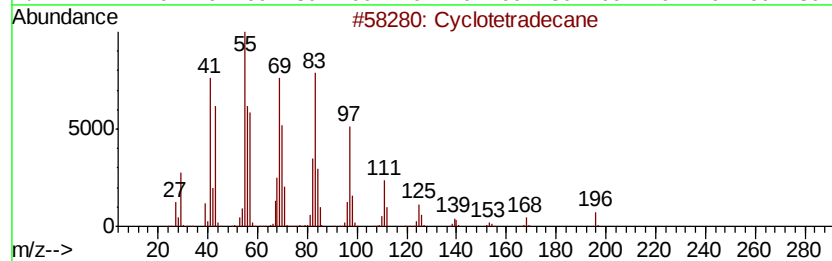
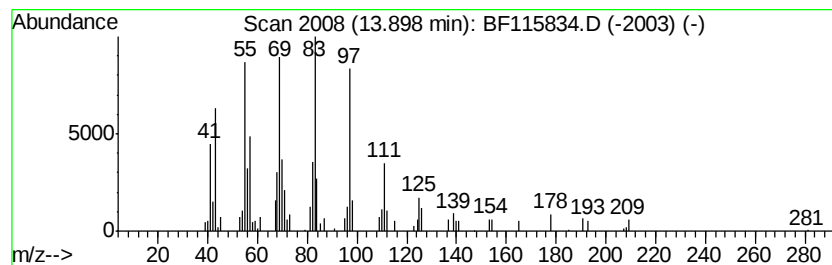
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Peak Number 6 Cyclotetradecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.90	3.21 ng	195509	Chrysene-d12		14.02
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cvcclotetradecane	196	C14H28	000295-17-0	72
2	Cycloeicosane	280	C20H40	000296-56-0	58
3	1-Hexadecanol	242	C16H34O	036653-82-4	58
4	Cvclopentane, undecyl-	224	C16H32	006785-23-5	49
5	Acetic acid, trifluoro-, undecyl...	268	C13H23F3O2	053800-01-4	46



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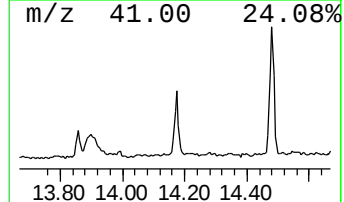
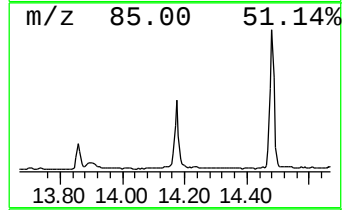
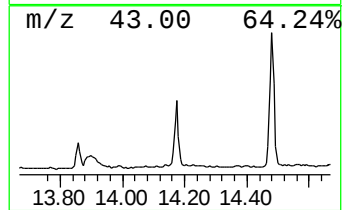
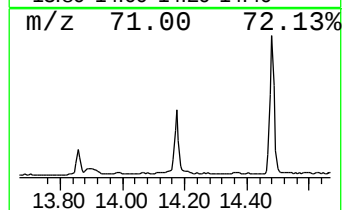
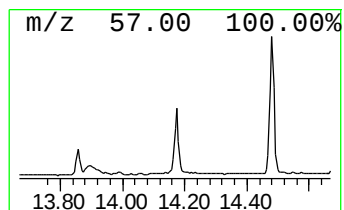
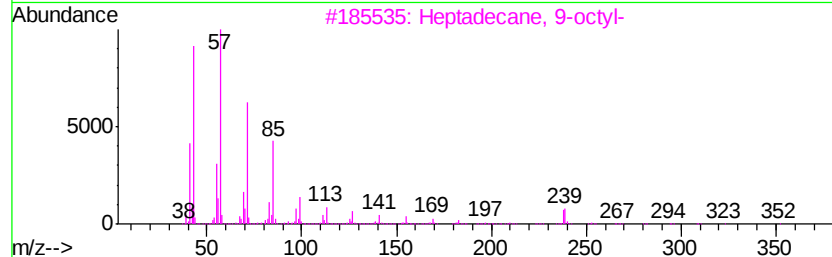
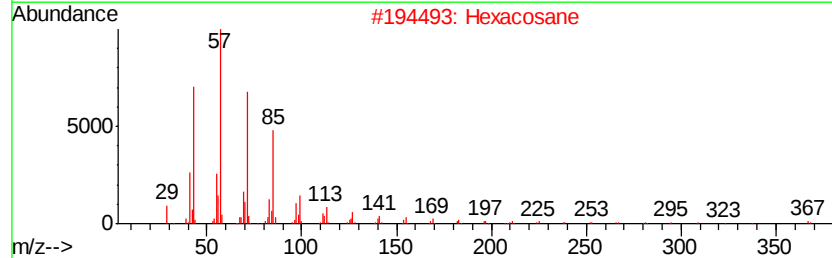
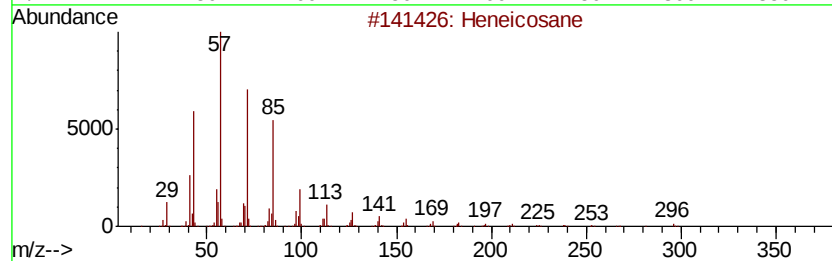
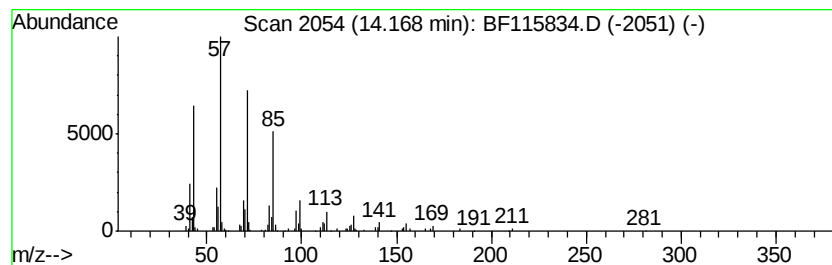
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 7 Heneicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.17	3.59 ng	218707	Chrysene-d12	14.02

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Hexacosane		366	C26H54	000630-01-3	91
3	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
4	Octacosane		394	C28H58	000630-02-4	91
5	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072919\
Data File : BF115834.D
Acq On : 29 Jul 2019 16:14
Operator : HP/JU
Sample : K4036-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

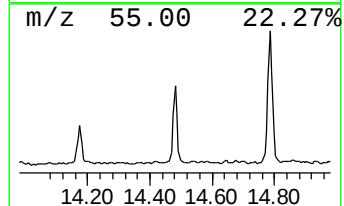
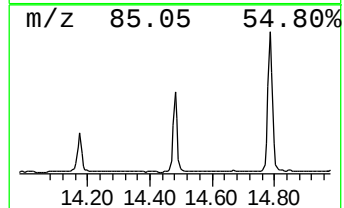
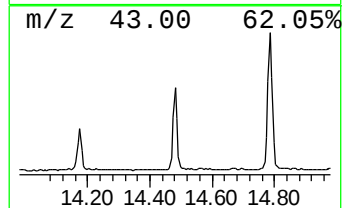
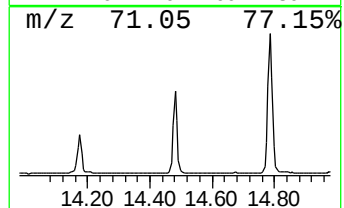
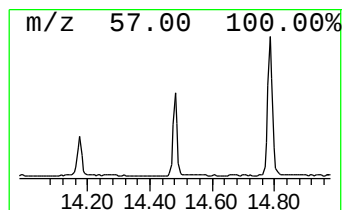
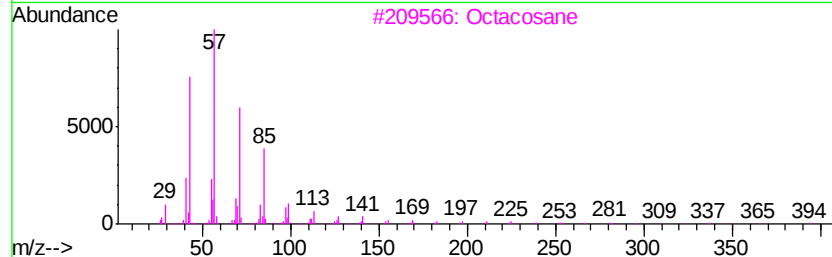
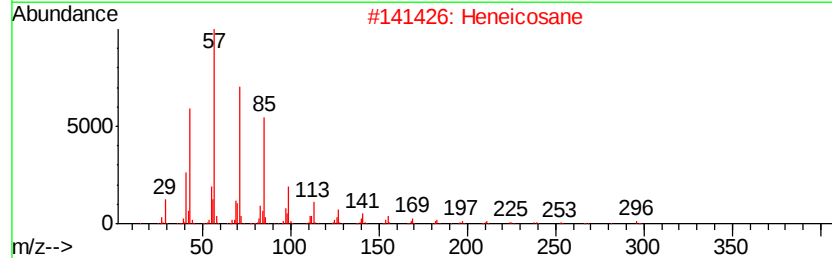
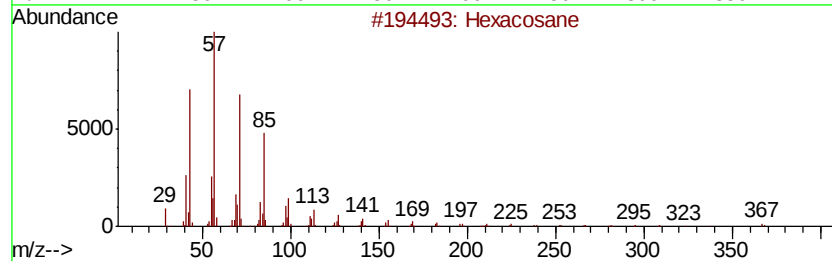
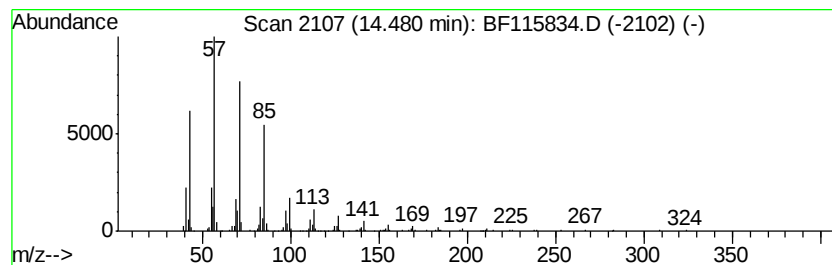
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.48	7.76 ng	472292	Chrysene-d12	14.02
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexacosane	366 C26H54	000630-01-3	91
2	Heneicosane	296 C21H44	000629-94-7	91
3	Octacosane	394 C28H58	000630-02-4	91
4	Heptadecane, 9-octyl-	352 C25H52	007225-64-1	91
5	Docosane, 11-butyl-	366 C26H54	013475-76-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072919\
Data File : BF115834.D
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Misc :
ALS Vial : 4 Sample Multiplier: 1

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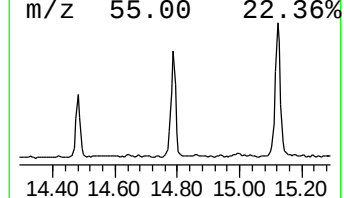
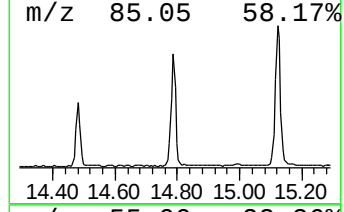
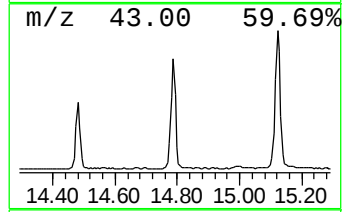
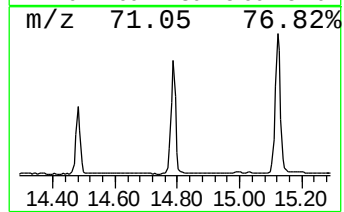
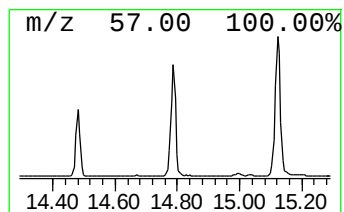
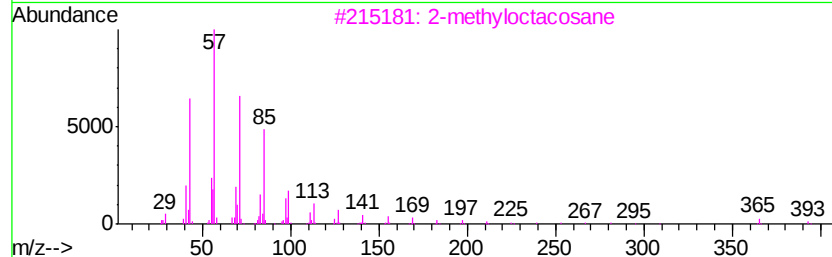
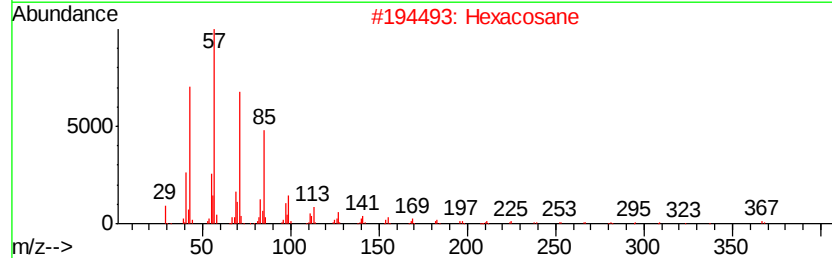
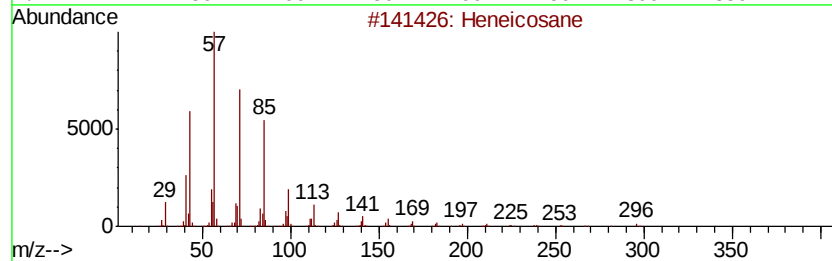
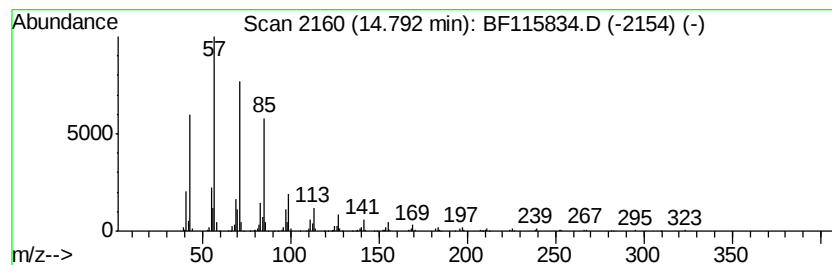
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane, 7-hexyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.79	6.44 ng	859694	Perylene-d12	15.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5		2-Bromotetradecane	276	C14H29Br	074036-95-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072919\
Data File : BF115834.D
Acq On : 29 Jul 2019 16:14
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Instrument :
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ClientSampleId :
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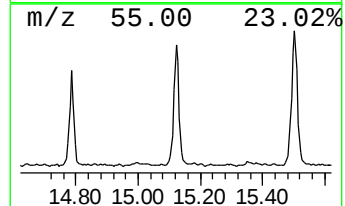
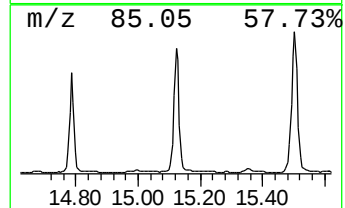
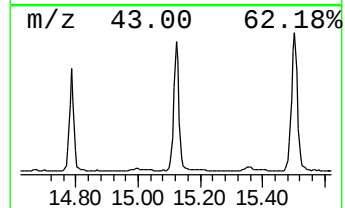
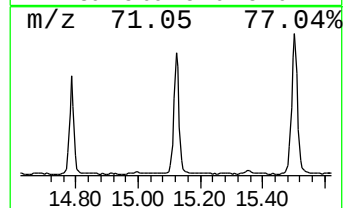
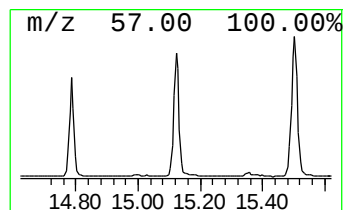
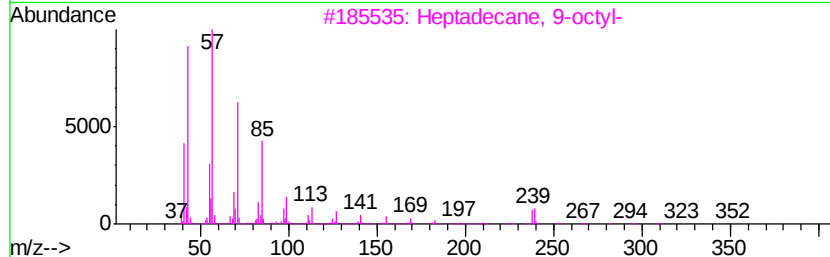
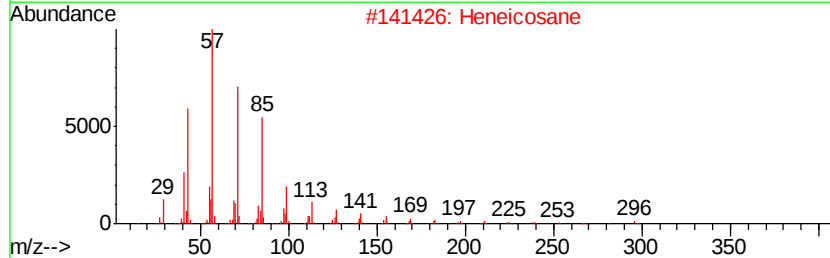
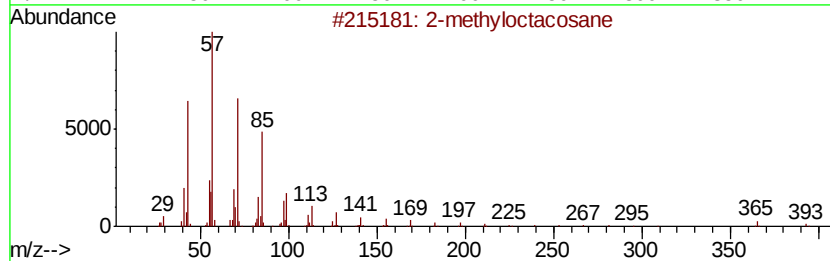
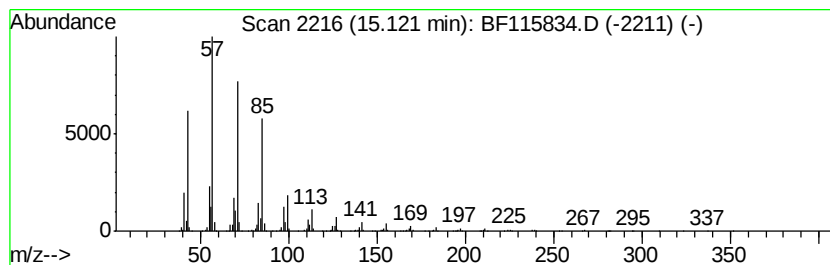
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 2-methyloctacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.12	9.44 ng	1259260	Perylene-d12	15.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		Octadecane	254	C18H38	000593-45-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072919\
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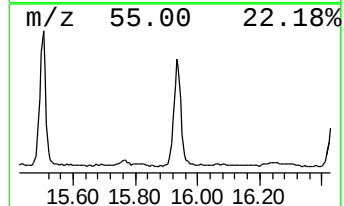
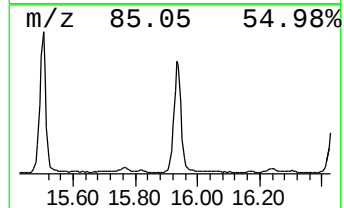
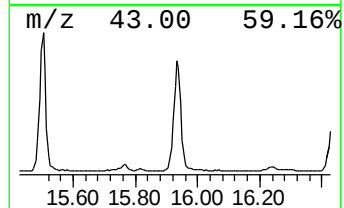
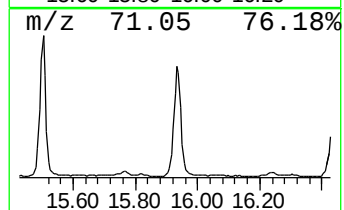
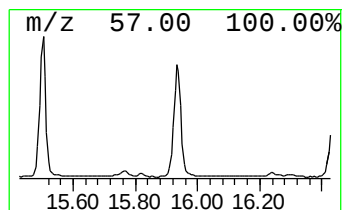
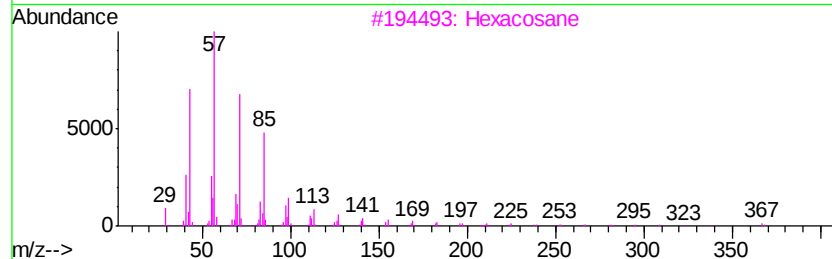
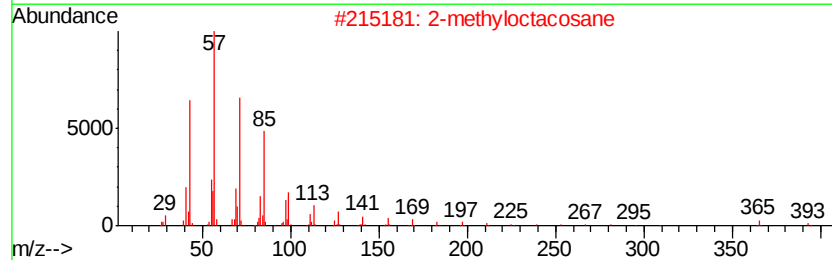
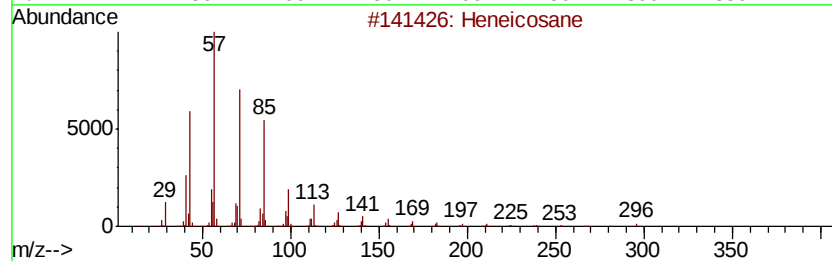
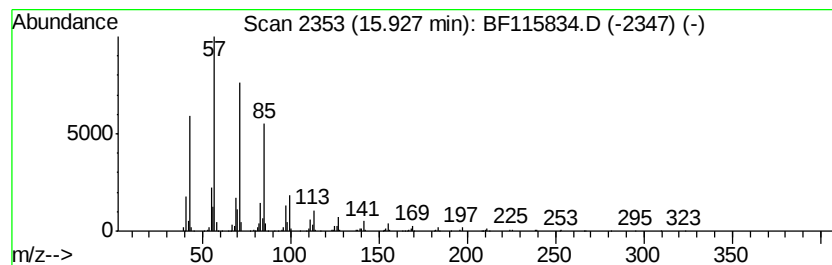
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.93	11.85 ng	1581490	Perylene-d12	15.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072919\
Data File : BF115834.D
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Sample : K4036-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

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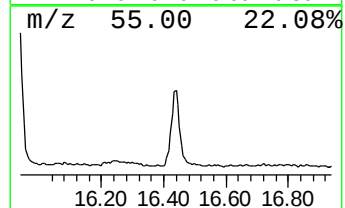
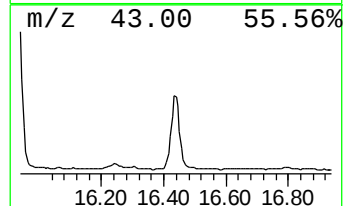
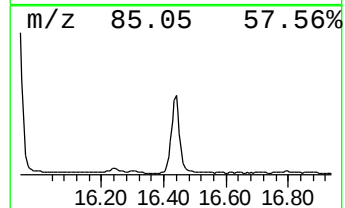
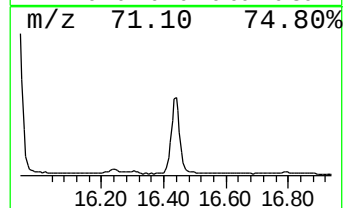
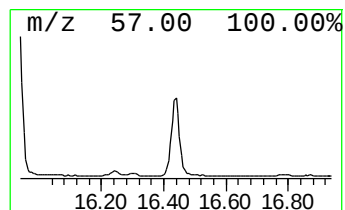
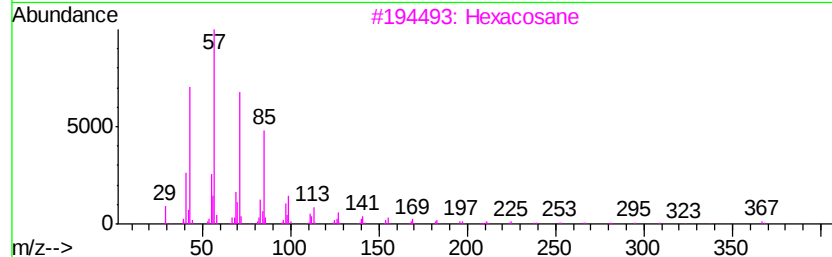
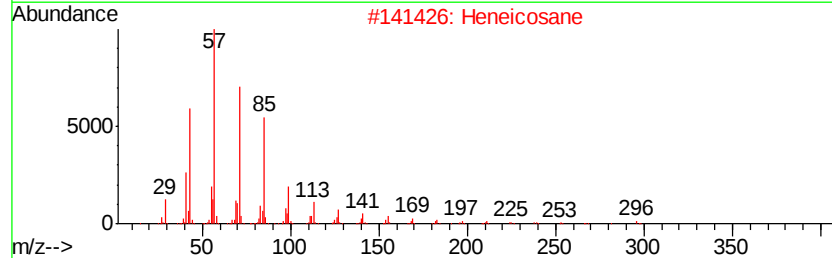
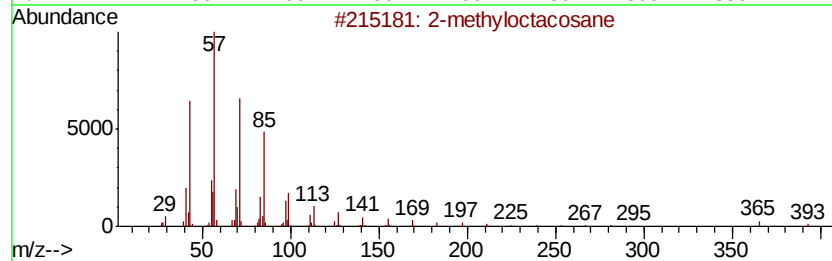
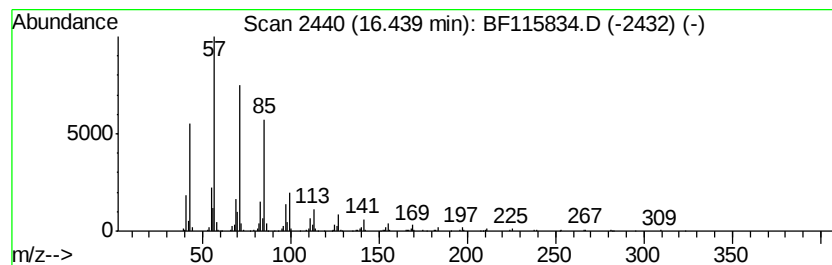
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptadecane, 9-octyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.44	7.29 ng	972062	Perylene-d12	15.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072919\
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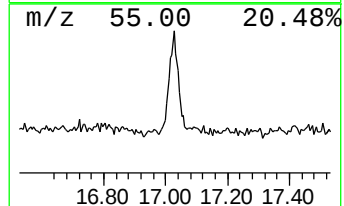
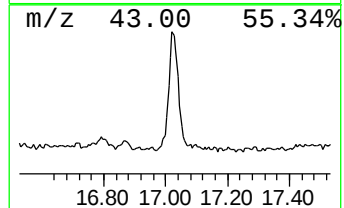
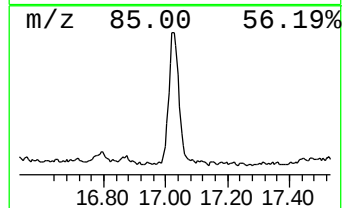
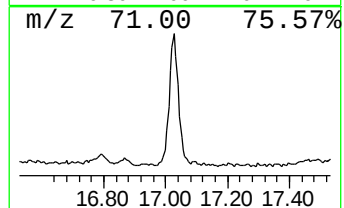
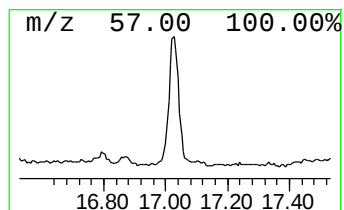
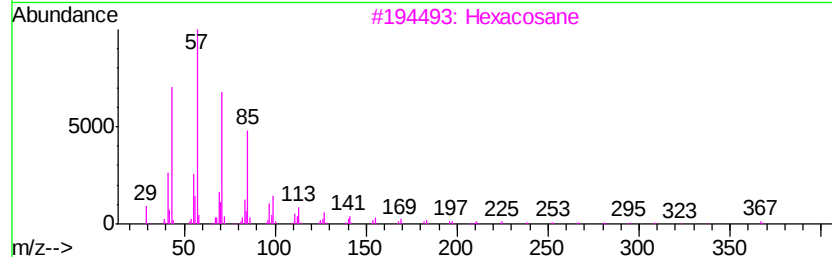
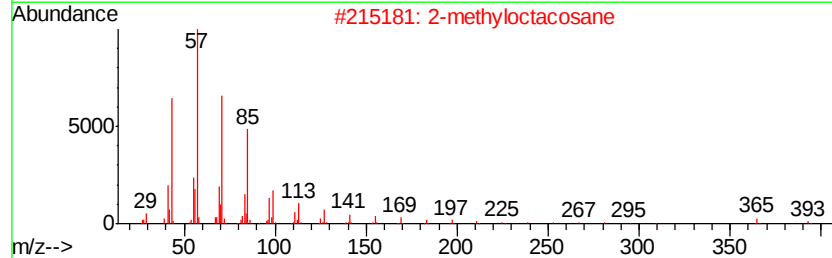
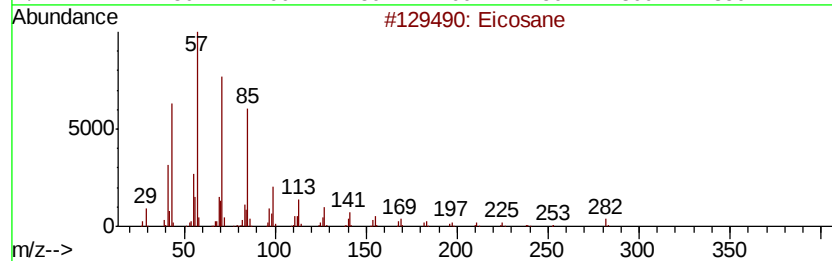
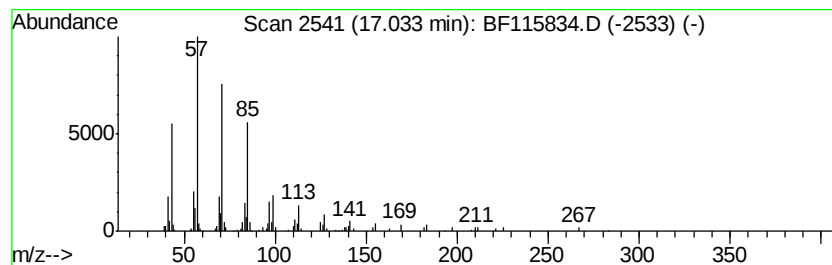
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Eicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.03	2.12 ng	283061	Perylene-d12	15.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072919\
Data File : BF115834.D
Acq On : 29 Jul 2019 16:14
Operator : HP/JU
Sample : K4036-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WS-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF071219.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
Butane, 2-methoxy...	2.13	81.9	ng	3356110	1	6.86	819387	20.0
unknown6.63	6.63	76.5	ng	3134630	1	6.86	819387	20.0
n-Hexadecanoic acid	11.90	6.1	ng	395839	4	11.37	1292600	20.0
Octadecanoic acid	12.69	10.9	ng	702571	4	11.37	1292600	20.0
Cyclotetradecane	13.90	3.2	ng	195509	5	14.02	1216860	20.0
Heneicosane	14.17	3.6	ng	218707	5	14.02	1216860	20.0
Hexacosane	14.48	7.8	ng	472292	5	14.02	1216860	20.0
Eicosane, 7-hexyl-	14.79	6.4	ng	859694	6	15.48	2668470	20.0
2-methyloctacosane	15.12	9.4	ng	1259260	6	15.48	2668470	20.0
Heptacosane	15.93	11.9	ng	1581490	6	15.48	2668470	20.0
Heptadecane, 9-oc...	16.44	7.3	ng	972062	6	15.48	2668470	20.0
Eicosane	17.03	2.1	ng	283061	6	15.48	2668470	20.0