

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF012020\
 Data File : BF118595.D
 Acq On : 20 Jan 2020 17:07
 Operator : CG/JU
 Sample : L1154-07
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B3-15-20

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-BF012020.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	2.216	16	22	53	rVB	2481341	4037389	7.22%	1.383%
2	5.098	506	512	521	rBV	1261040	1490471	2.67%	0.511%
3	5.498	575	580	599	rBV	5280552	9144579	16.35%	3.133%
4	5.628	599	602	622	rVB	1001624	1565167	2.80%	0.536%
5	6.504	745	751	755	rBV	9015142	10829142	19.37%	3.710%
6	6.651	771	776	779	rBV	10422477	10748872	19.22%	3.683%
7	6.881	811	815	818	rBV	3254055	2739376	4.90%	0.939%
8	7.040	837	842	845	rBV	9299046	8483380	15.17%	2.907%
9	7.440	904	910	913	rBV	6336316	6658935	11.91%	2.281%
10	8.163	1028	1033	1036	rBV	3766782	3501808	6.26%	1.200%
11	9.239	1210	1216	1219	rBV	12830073	12767619	22.83%	4.374%
12	9.628	1277	1282	1285	rBV	1065356	840023	1.50%	0.288%
13	9.916	1324	1331	1335	rBV	4865126	4372033	7.82%	1.498%
14	10.704	1459	1465	1468	rBV	9207060	9035725	16.16%	3.096%
15	11.404	1579	1584	1587	rBV	4816717	4645856	8.31%	1.592%
16	11.451	1587	1592	1593	rBV	498545	613804	1.10%	0.210%
17	11.545	1604	1608	1610	rBV	537281	694192	1.24%	0.238%
18	11.580	1610	1614	1616	rBV	1460395	2329315	4.17%	0.798%
19	11.775	1645	1647	1649	rBV	1202707	983866	1.76%	0.337%
20	11.933	1668	1674	1683	rVB	1676337	2077812	3.72%	0.712%
21	12.716	1801	1807	1819	rVB	1045851	1233777	2.21%	0.423%
22	12.927	1837	1843	1849	rVV	721733	2285821	4.09%	0.783%
23	12.992	1849	1854	1857	rVB	13054361	13464898	24.08%	4.613%
24	13.092	1857	1871	1885	rVB	499398	1881457	3.36%	0.645%
25	13.463	1909	1934	1946	rBV	13648214	55913197	100.00%	19.157%
26	13.569	1948	1952	1956	rVB	774844	831161	1.49%	0.285%
27	13.886	2002	2006	2011	rVV2	2938492	4293038	7.68%	1.471%
28	13.980	2013	2022	2028	rVV	1209177	3785563	6.77%	1.297%
29	14.039	2028	2032	2035	rVV	3878251	4190691	7.49%	1.436%
30	14.080	2035	2039	2050	rVB	572731	1299449	2.32%	0.445%
31	14.210	2056	2061	2064	rBV	4472098	5138411	9.19%	1.760%
32	14.292	2064	2075	2092	rVB	11342383	28555285	51.07%	9.783%
33	14.516	2108	2113	2117	rBV	7673143	8815108	15.77%	3.020%
34	14.827	2160	2166	2175	rVB	9007770	11612156	20.77%	3.978%

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Integration Parameters: rteint.p
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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

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Peak separation: 5

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35	14.951	2177	2187	2196	rVB	3517410	8672696	15.51%	2.971%
36	15.174	2217	2225	2235	rVB	9178831	12486226	22.33%	4.278%
37	15.510	2277	2282	2285	rBV	2248936	2926189	5.23%	1.003%
38	15.563	2285	2291	2306	rVB	7177395	11135909	19.92%	3.815%
39	15.710	2309	2316	2327	rVB	806353	2121991	3.80%	0.727%
40	16.004	2357	2366	2371	rBV	4744161	8222100	14.71%	2.817%
41	16.515	2444	2453	2460	rBV	2157712	4279295	7.65%	1.466%
42	17.115	2548	2555	2563	rVB	547010	1171116	2.09%	0.401%

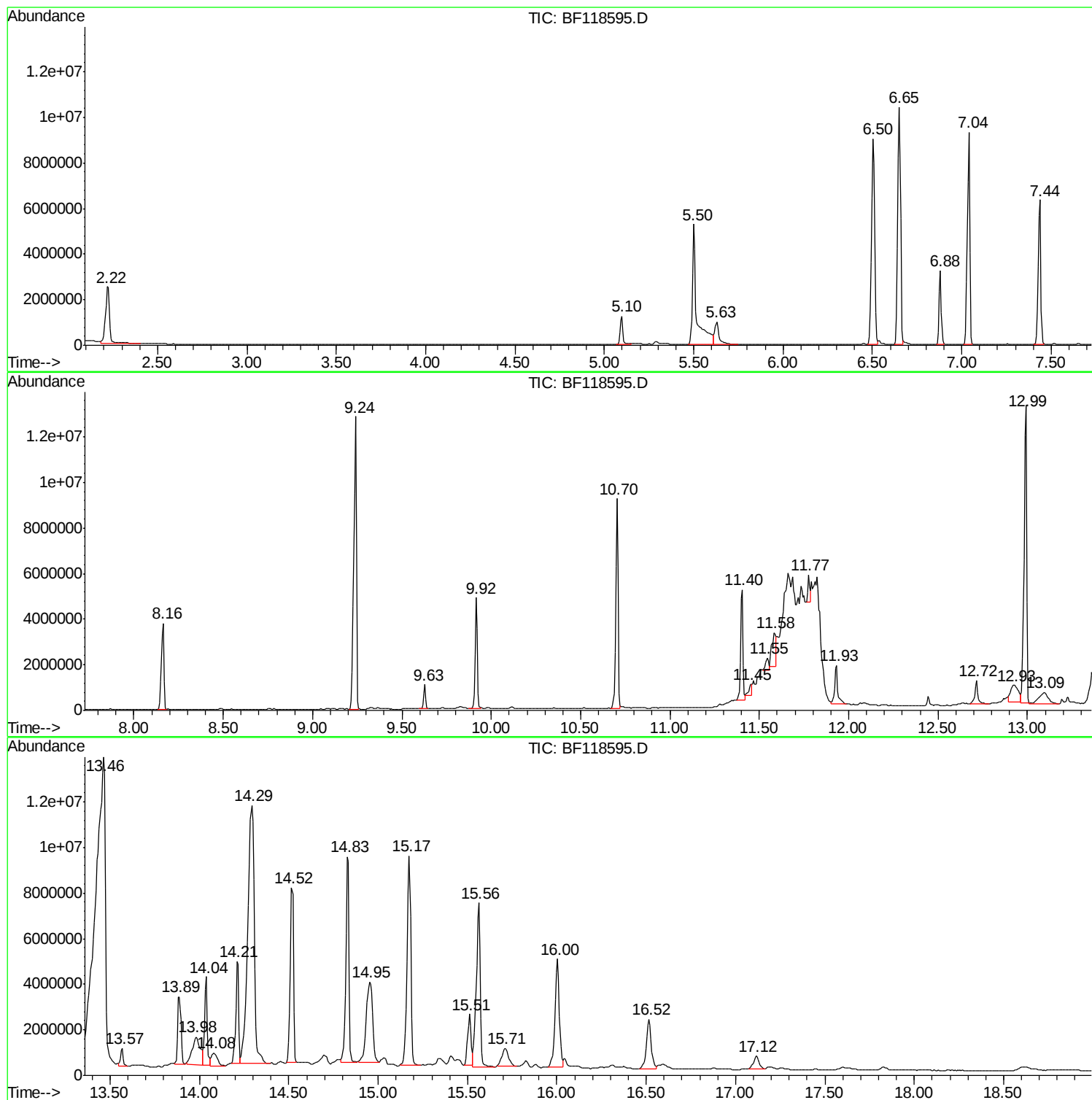
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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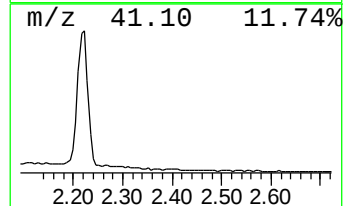
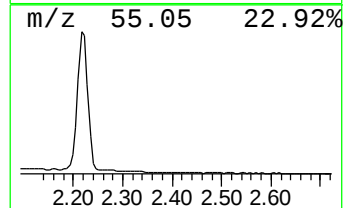
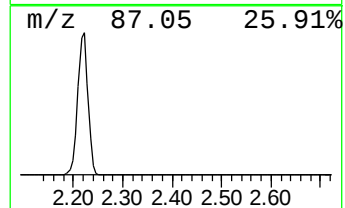
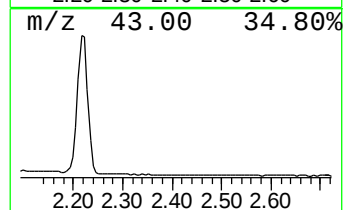
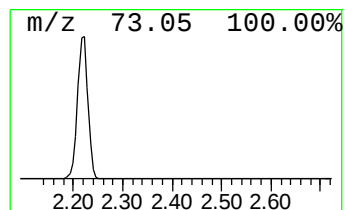
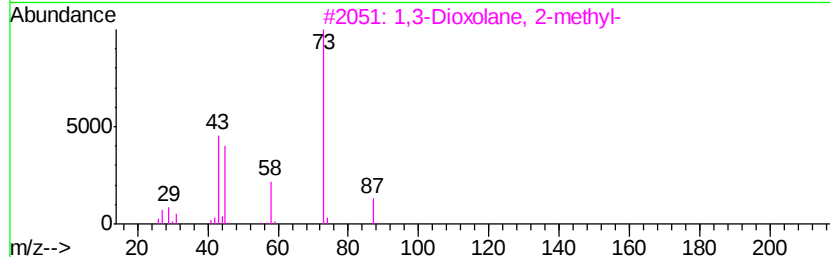
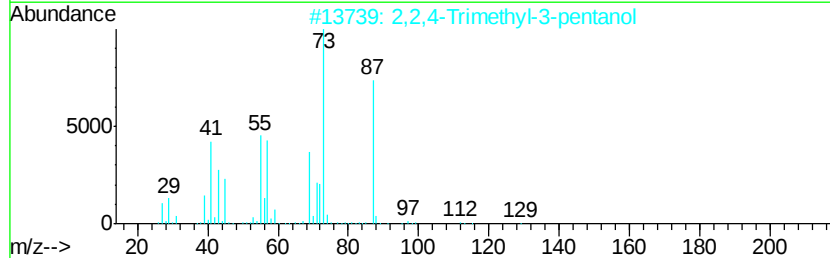
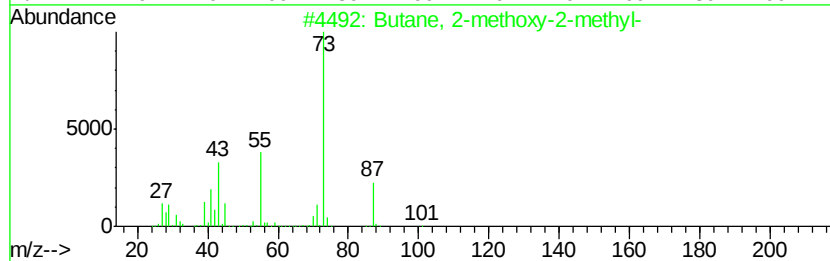
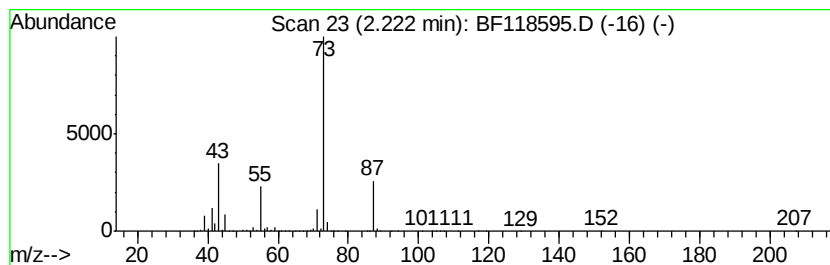
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.22	29.48 ng	4037390	1,4-Dichlorobenzene-d4	6.88

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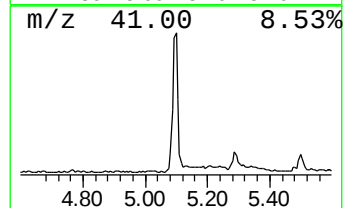
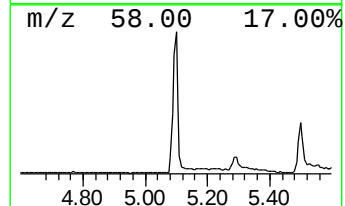
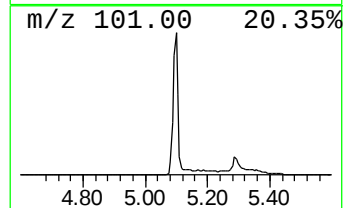
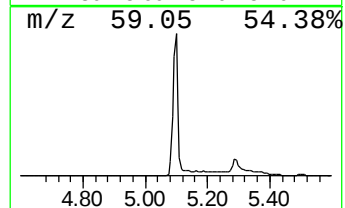
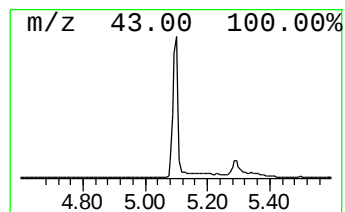
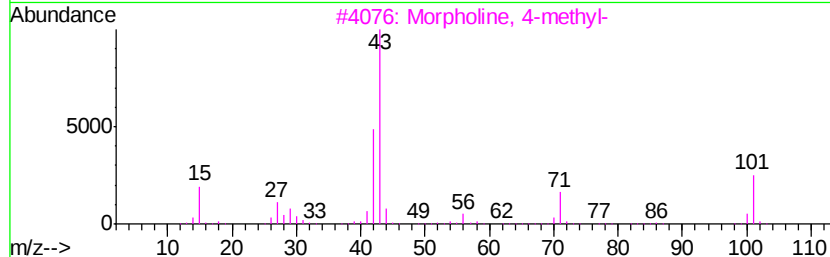
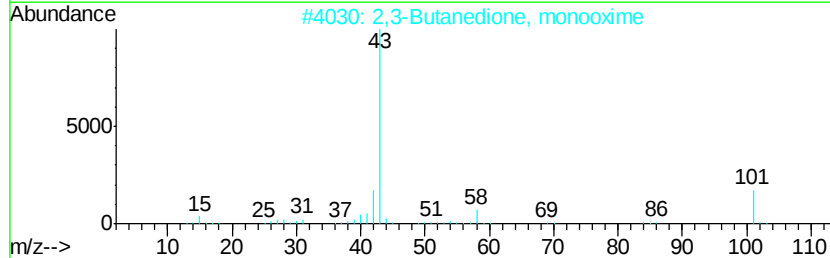
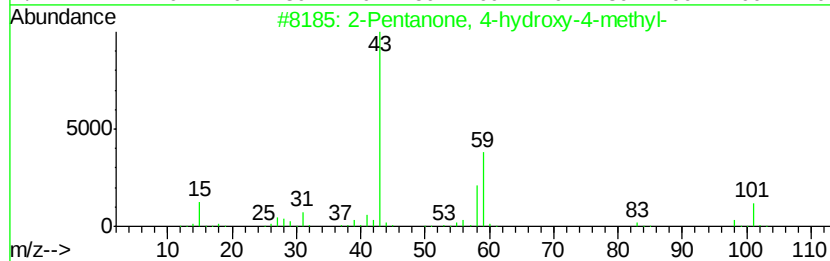
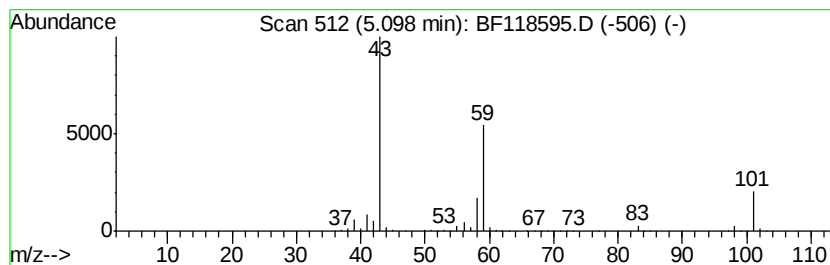
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	10.88 ng	1490470	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9



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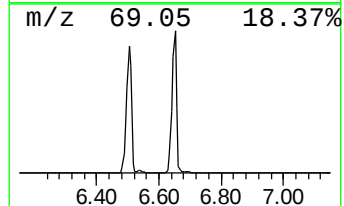
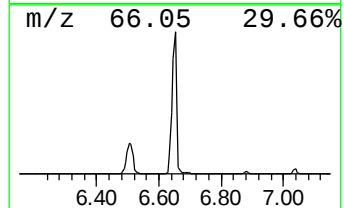
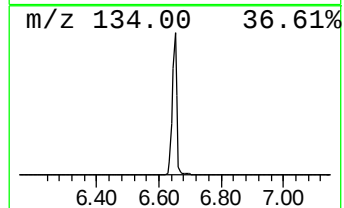
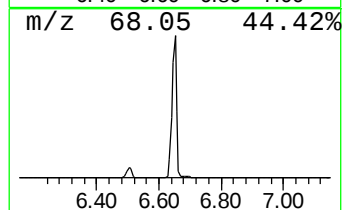
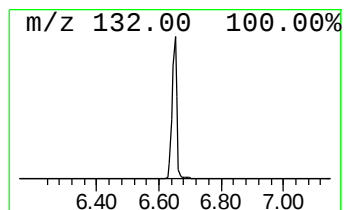
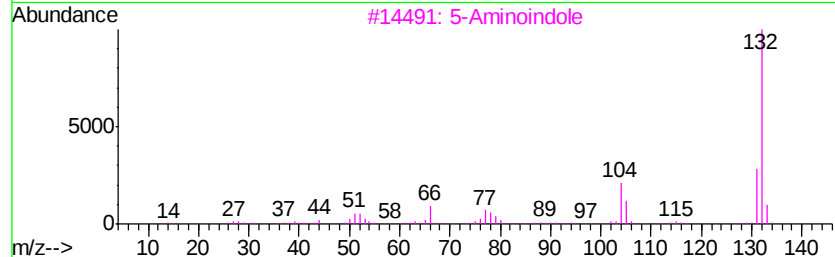
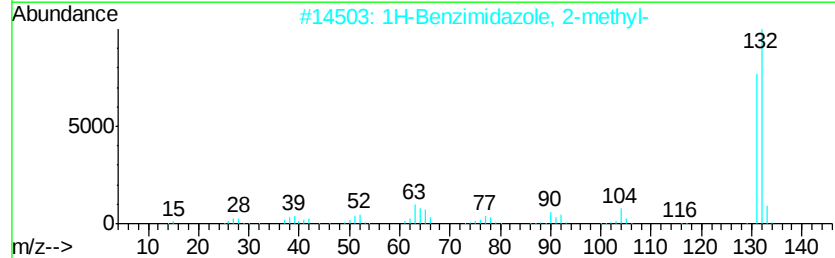
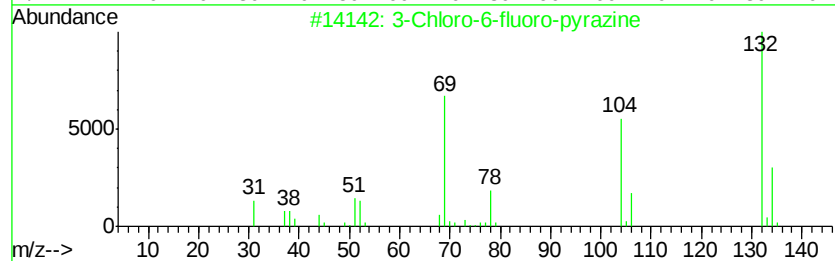
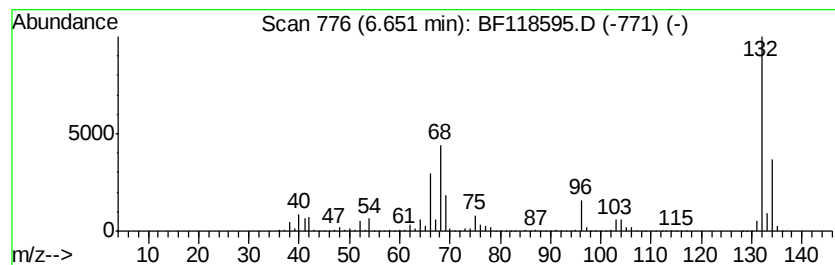
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.65 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	78.48 ng	10748900	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Aminoindole	132	C8H8N2	005192-03-0	22
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12



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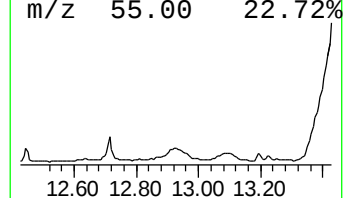
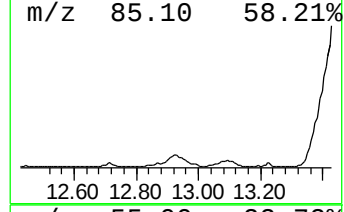
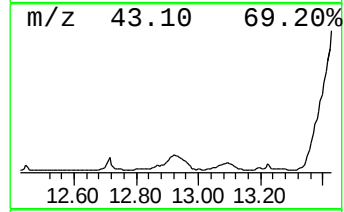
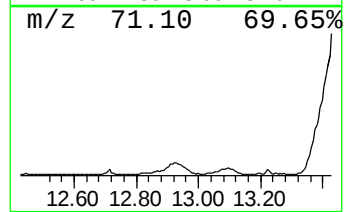
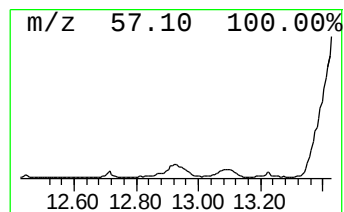
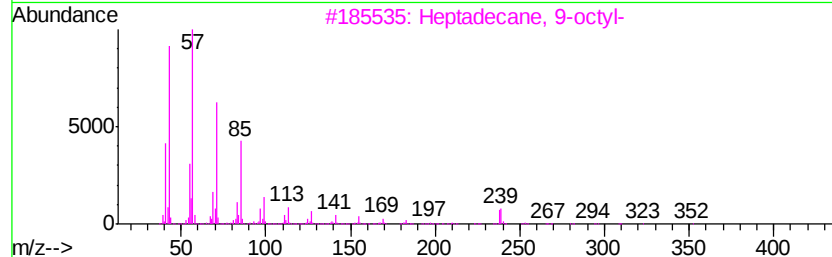
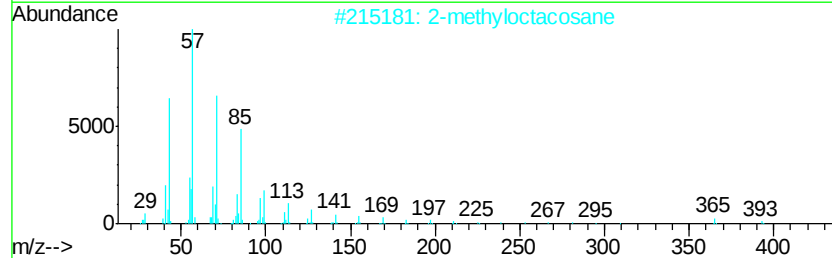
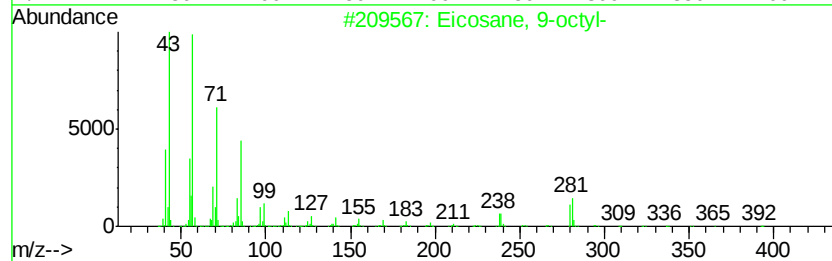
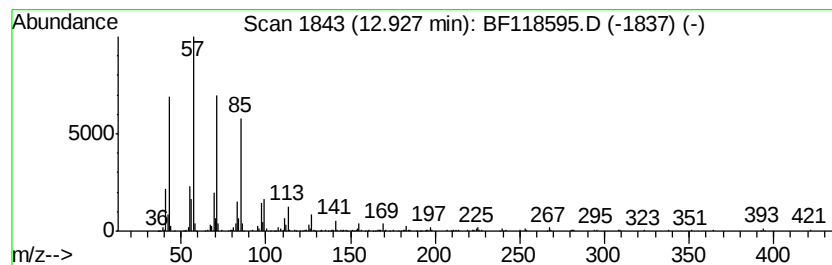
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Eicosane, 9-octyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.93	10.91 ng	2285820	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Tetracosane	338	C24H50	000646-31-1	91
5		Eicosane	282	C20H42	000112-95-8	90



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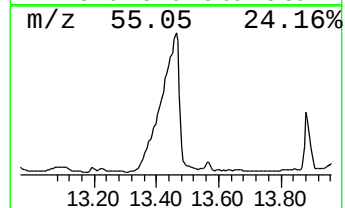
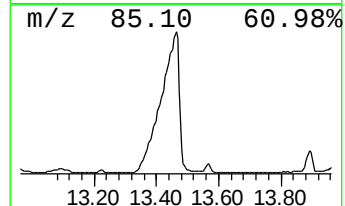
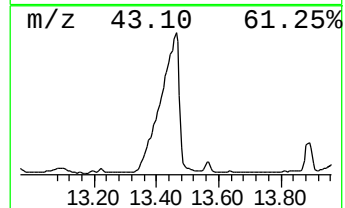
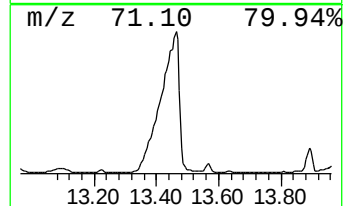
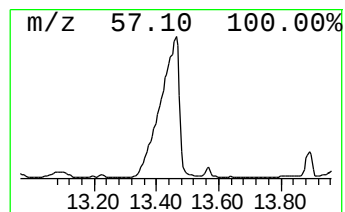
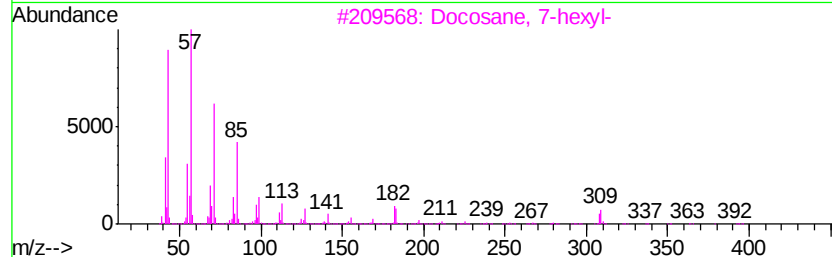
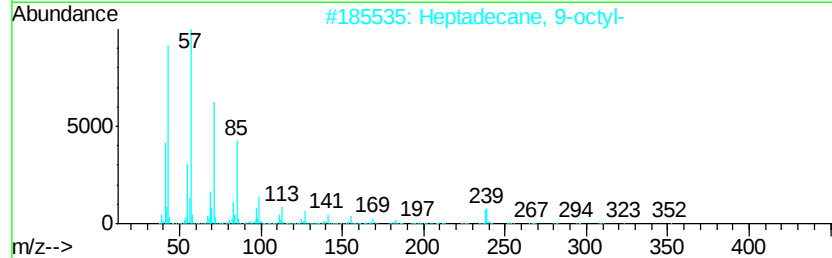
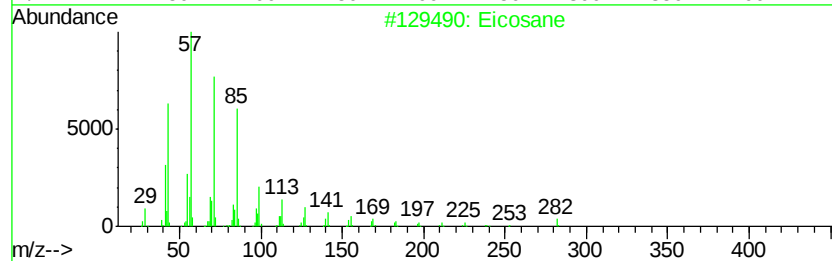
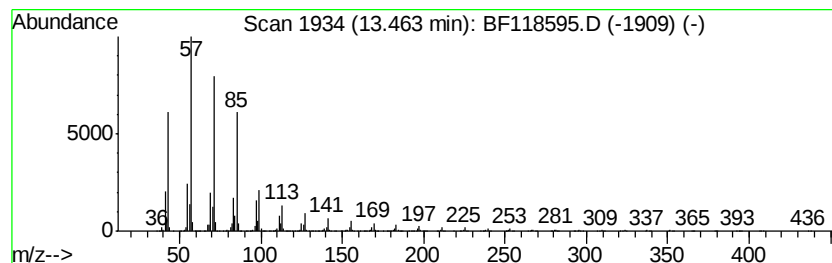
Instrument :
BNA_F
ClientSampleId :
B3-15-20

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.46	266.85 ng	55913200	Chrysene-d12	14.04
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	94
2	Heptadecane, 9-octyl-	352 C25H52	007225-64-1	93
3	Docosane, 7-hexyl-	394 C28H58	055373-86-9	93
4	Heneicosane	296 C21H44	000629-94-7	91
5	1-Iodo-2-methylundecane	296 C12H25I	073105-67-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118595.D
Acq On : 20 Jan 2020 17:07
Operator : CG/JU
Sample : L1154-07
Misc :
ALS Vial : 5 Sample Multiplier: 1

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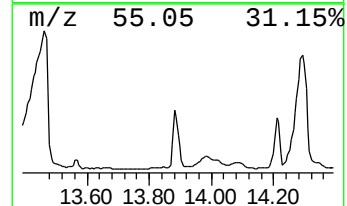
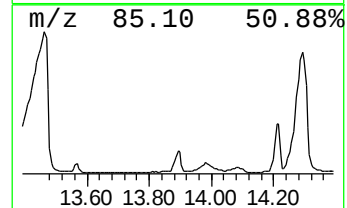
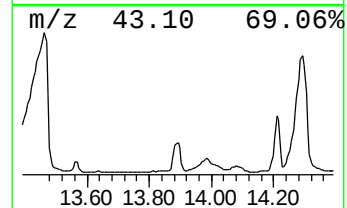
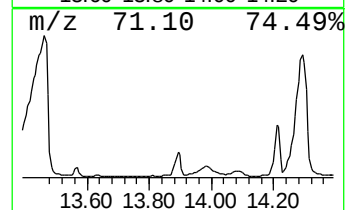
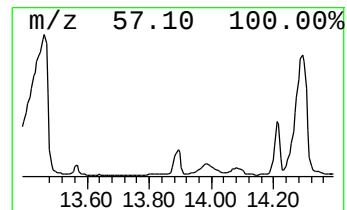
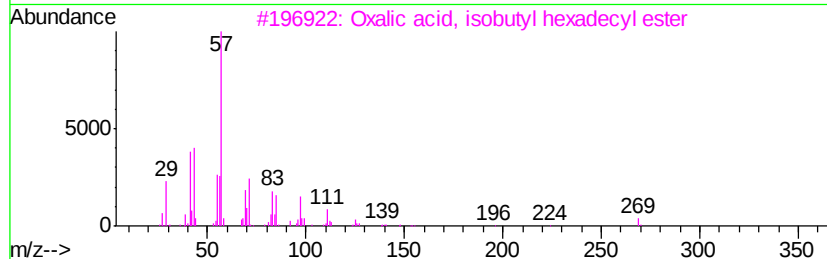
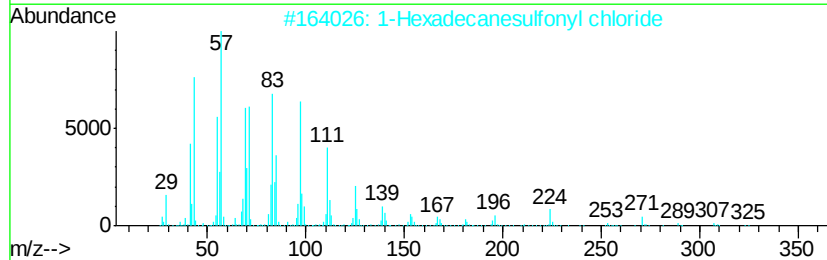
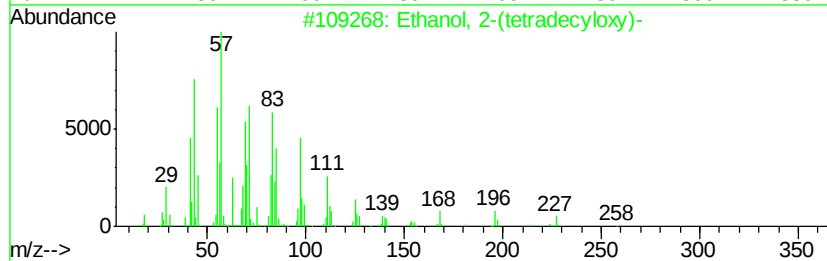
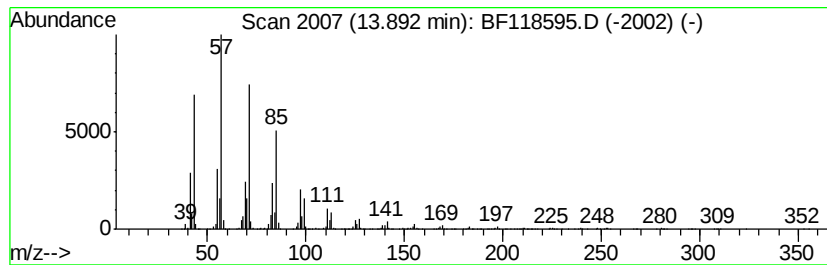
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Ethanol, 2-(tetradecyloxy)- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	20.49 ng	4293040	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
2		1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	91
3		Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
4		1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	90
5		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
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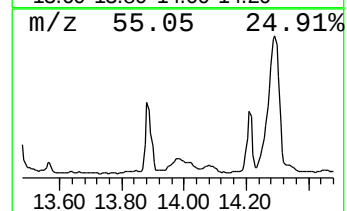
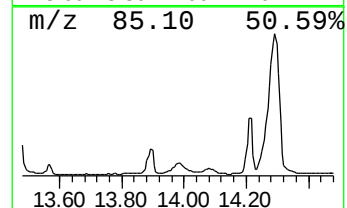
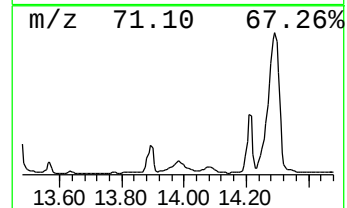
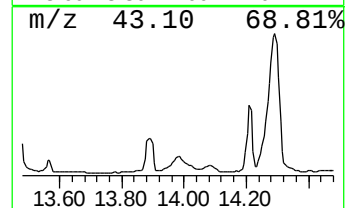
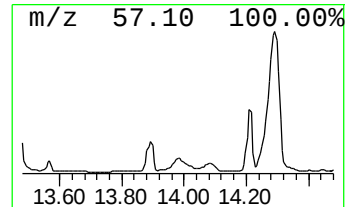
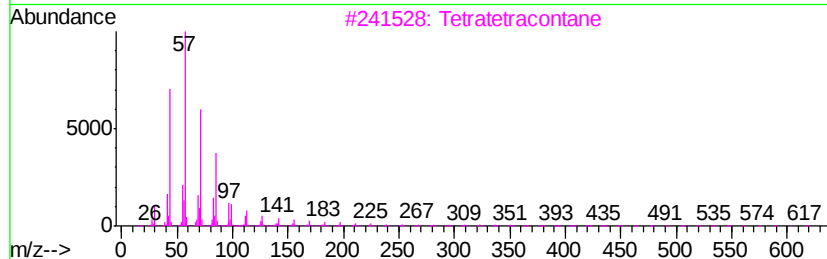
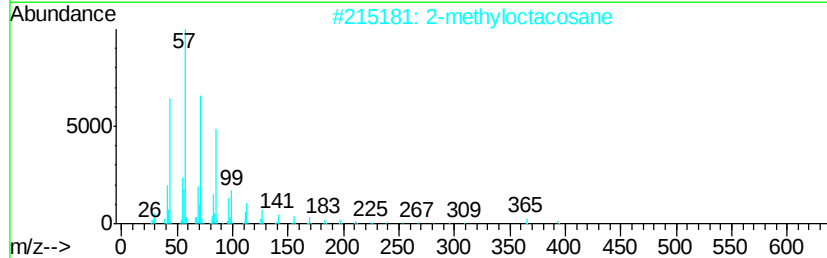
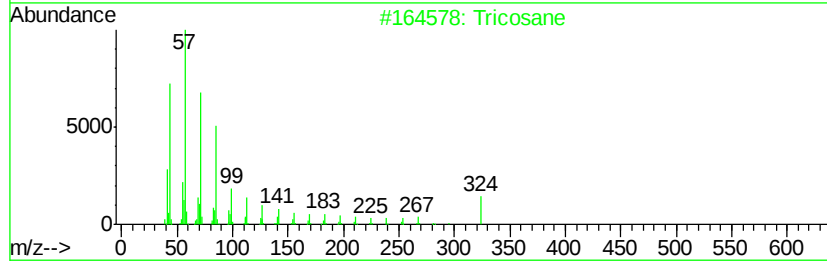
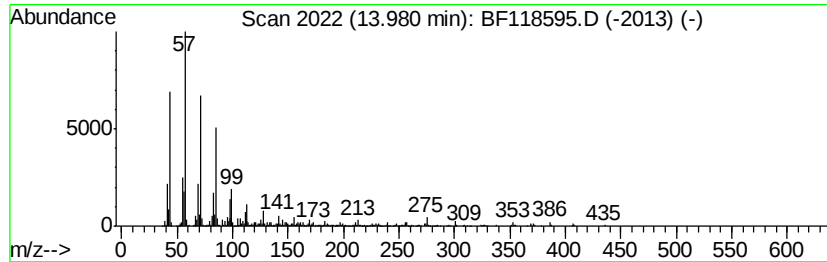
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tricosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.98	18.07 ng	3785560	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	97
2		2-methyloctacosane	408	C29H60	1000376-72-8	90
3		Tetratetracontane	619	C44H90	007098-22-8	90
4		Nonacosane	408	C29H60	000630-03-5	90
5		2-methylhexacosane	380	C27H56	1000376-72-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118595.D
Acq On : 20 Jan 2020 17:07
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Misc :
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ClientSampleId :
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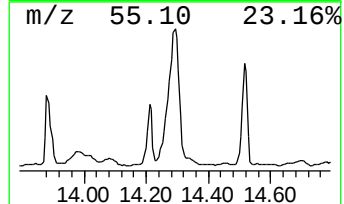
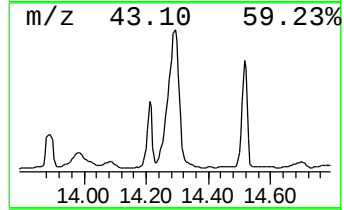
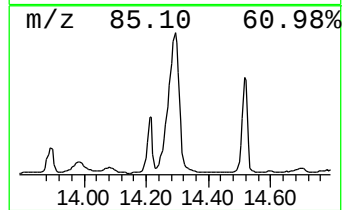
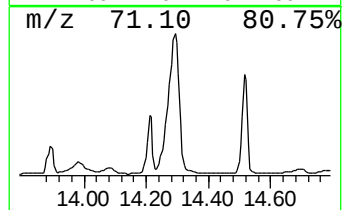
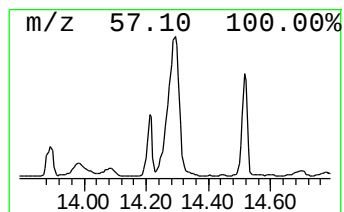
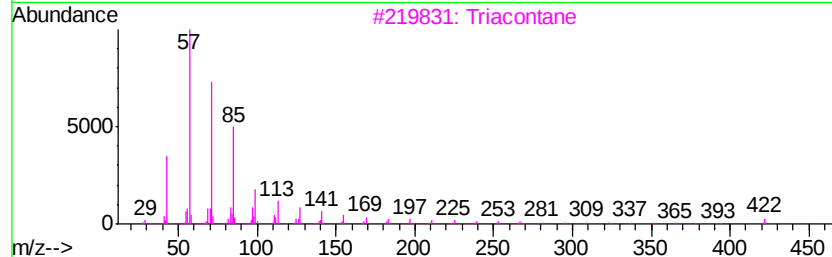
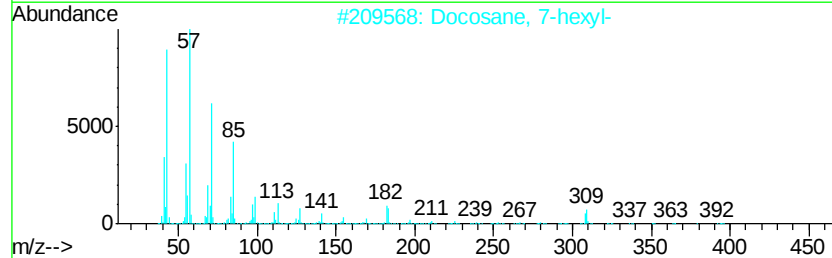
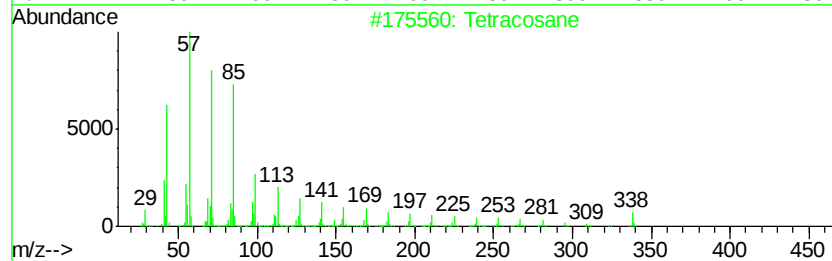
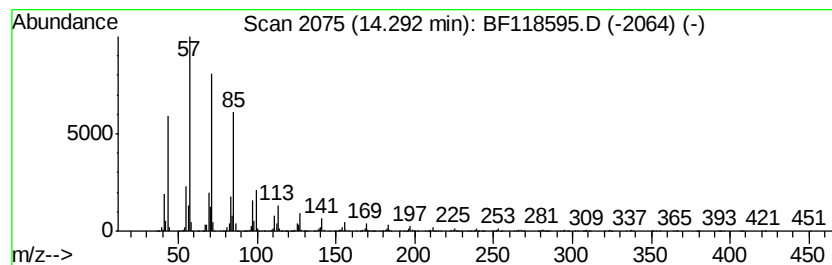
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetracosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	136.28 ng	28555300	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	95
2		Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
3		Triacontane	422	C30H62	000638-68-6	91
4		Tetratetracontane	619	C44H90	007098-22-8	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118595.D
Acq On : 20 Jan 2020 17:07
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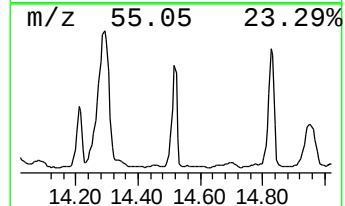
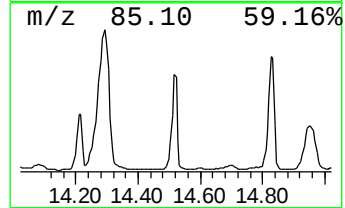
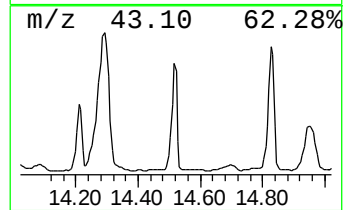
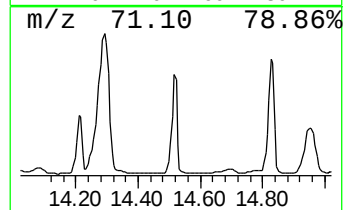
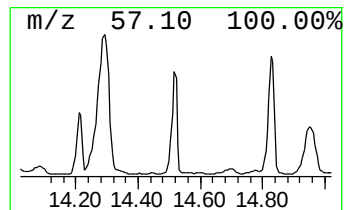
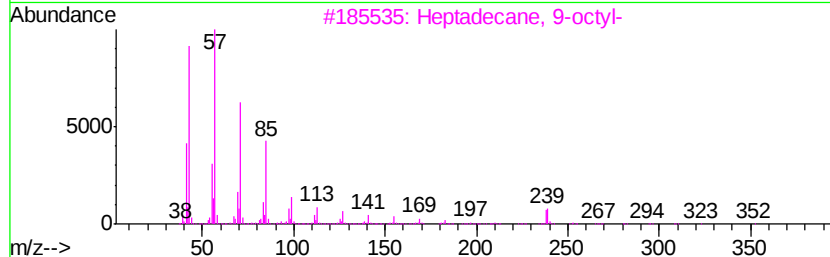
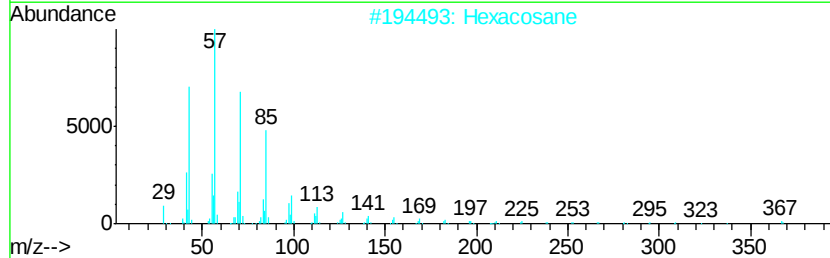
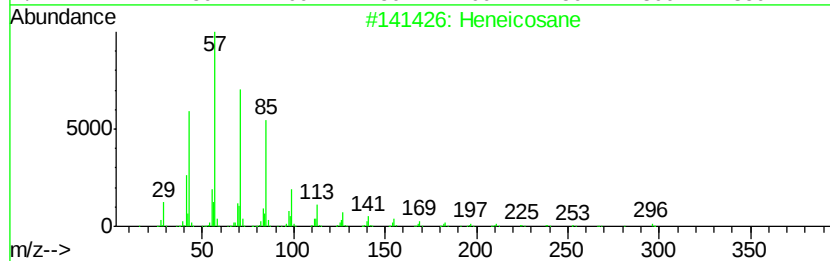
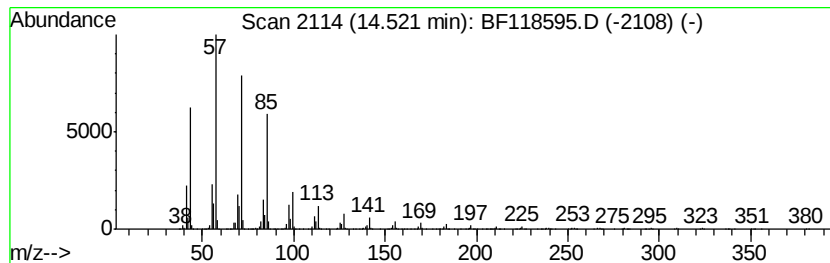
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	42.07 ng	8815110	Chrysene-d12	14.04

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		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118595.D
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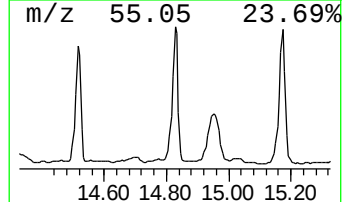
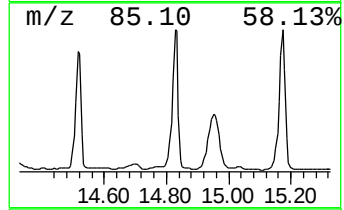
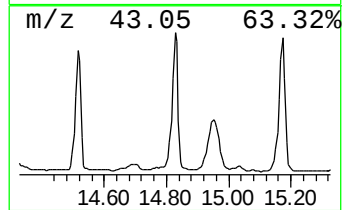
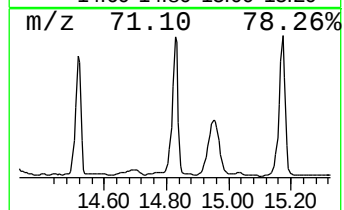
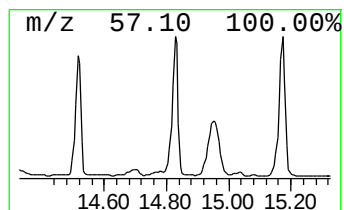
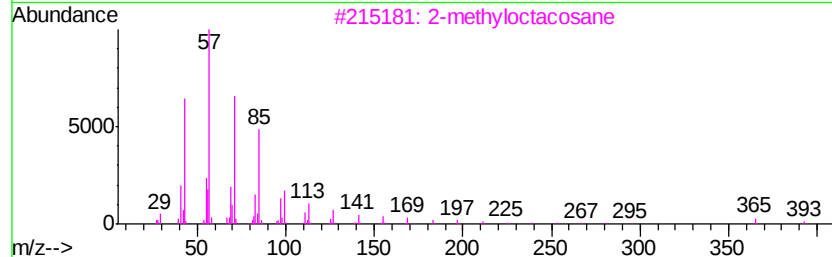
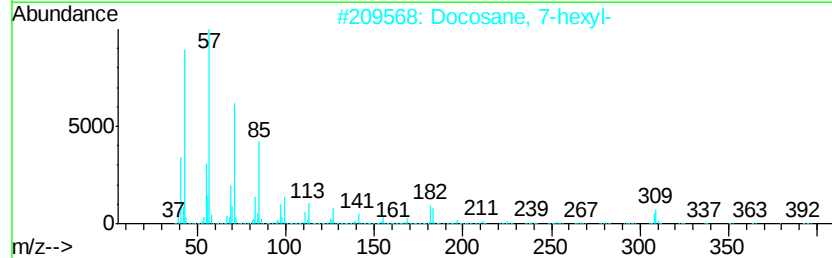
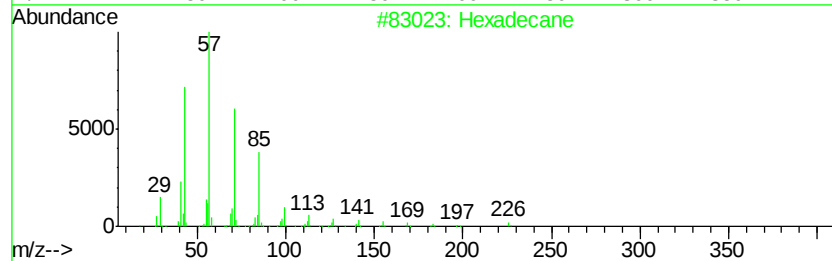
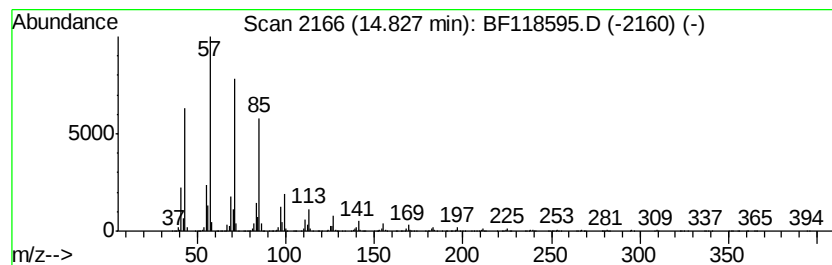
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Hexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	79.37 ng	11612200	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	95
2		Docosane, 7-hexyl-	394	C28H58	055373-86-9	94
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118595.D
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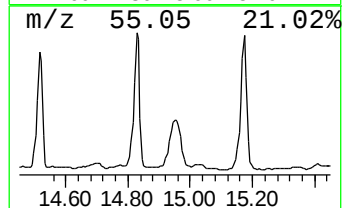
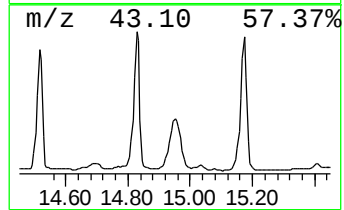
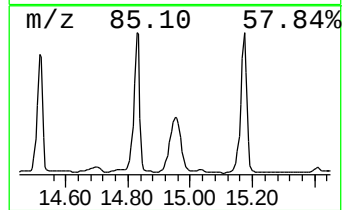
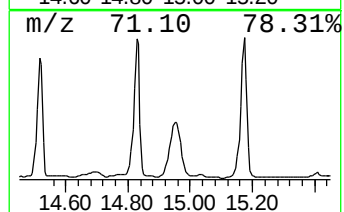
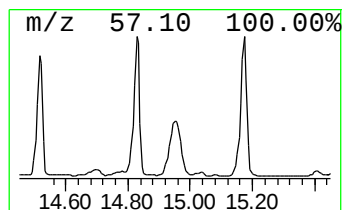
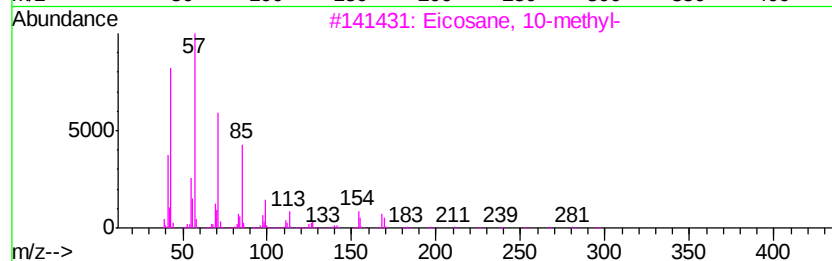
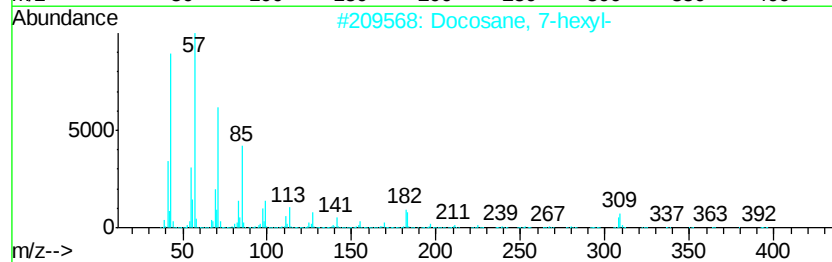
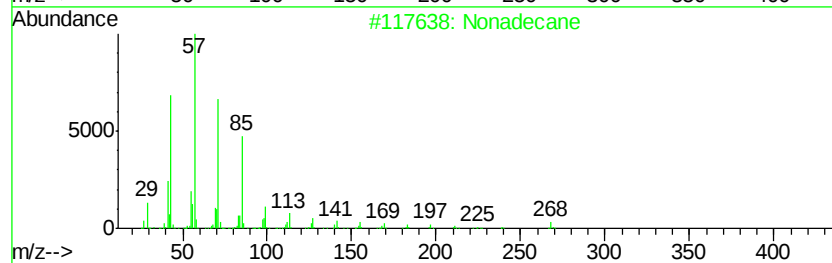
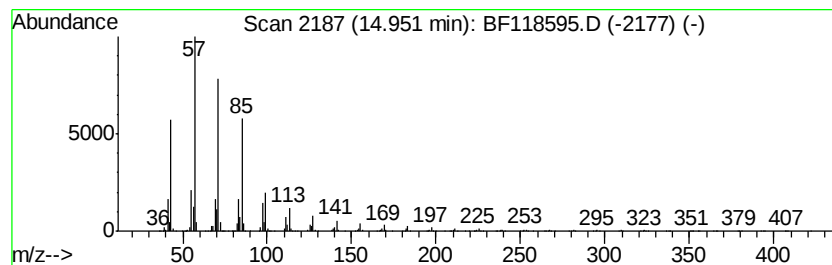
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Nonadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.95	59.28 ng	8672700	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	94
2		Docosane, 7-hexyl-	394	C28H58	055373-86-9	94
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Heneicosane	296	C21H44	000629-94-7	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
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Acq On : 20 Jan 2020 17:07
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ClientSampleId :
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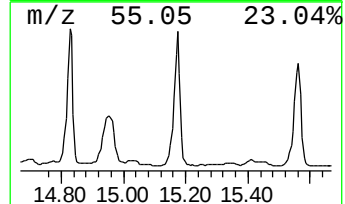
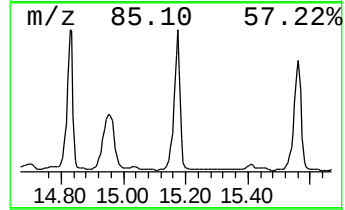
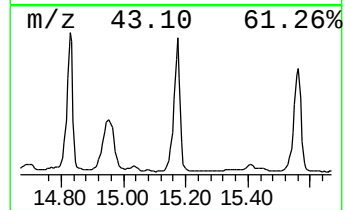
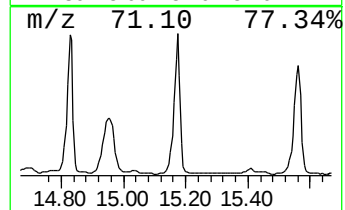
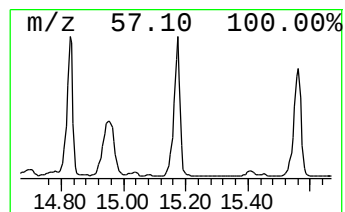
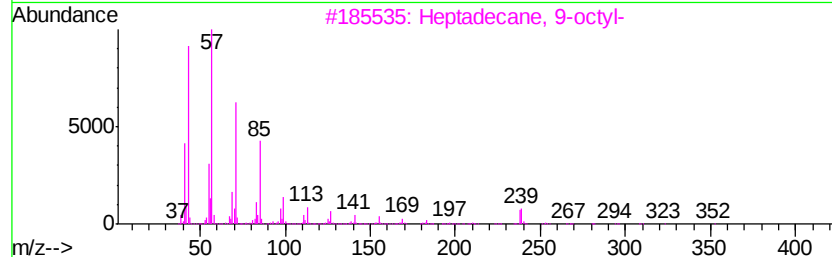
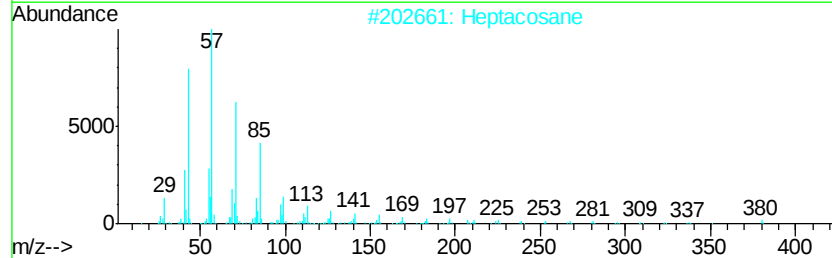
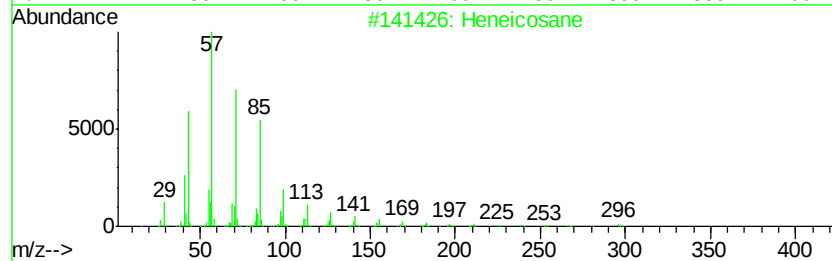
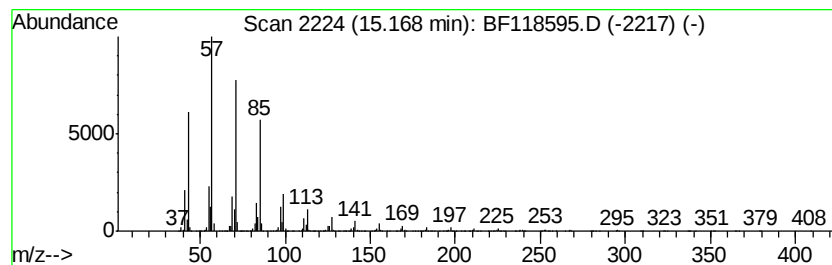
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Heptacosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	85.34 ng	12486200	Perylene-d12	15.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	91
2		Heptacosane	380	C27H56	000593-49-7	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118595.D
Acq On : 20 Jan 2020 17:07
Operator : CG/JU
Sample : L1154-07
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B3-15-20

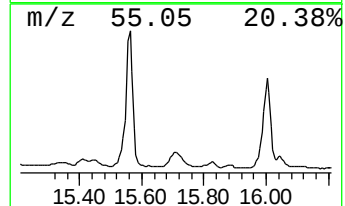
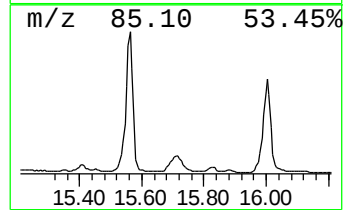
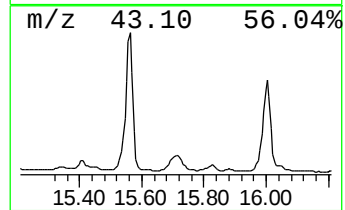
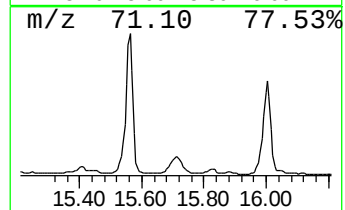
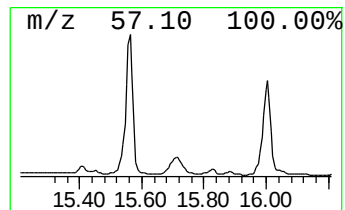
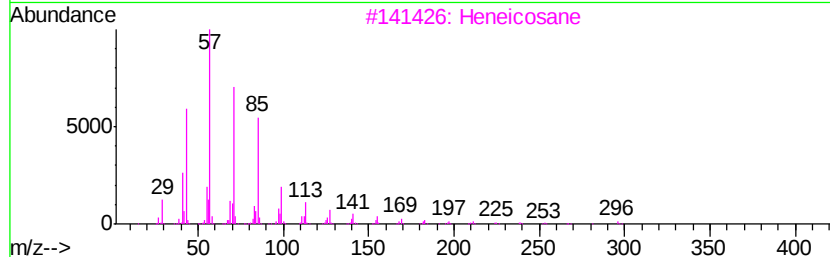
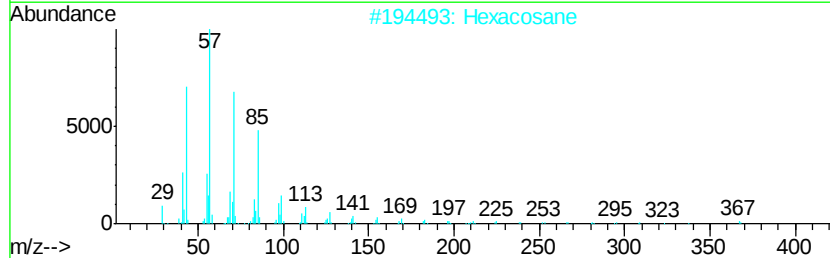
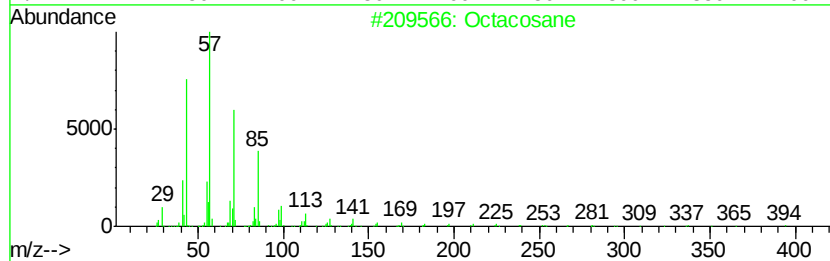
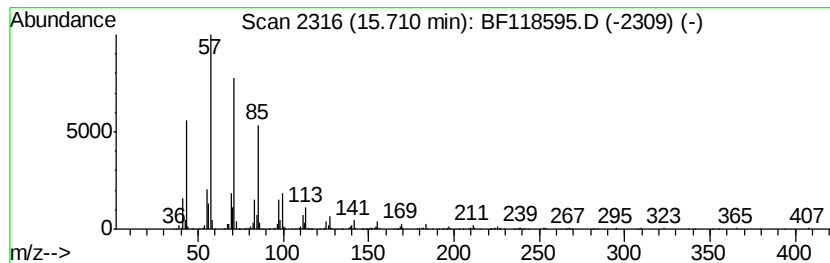
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Octacosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.71	14.50 ng	2121990	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118595.D
Acq On : 20 Jan 2020 17:07
Operator : CG/JU
Sample : L1154-07
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B3-15-20

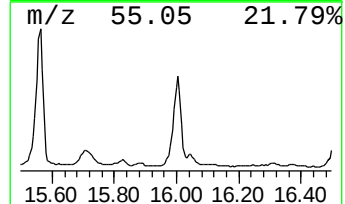
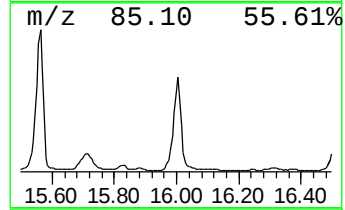
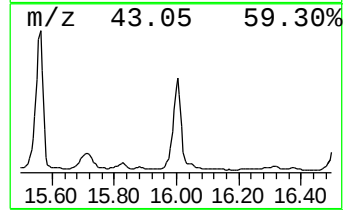
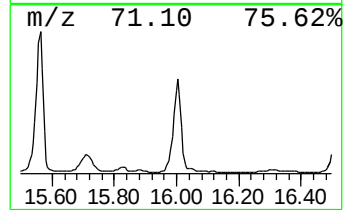
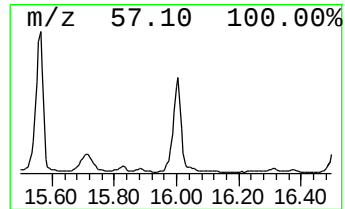
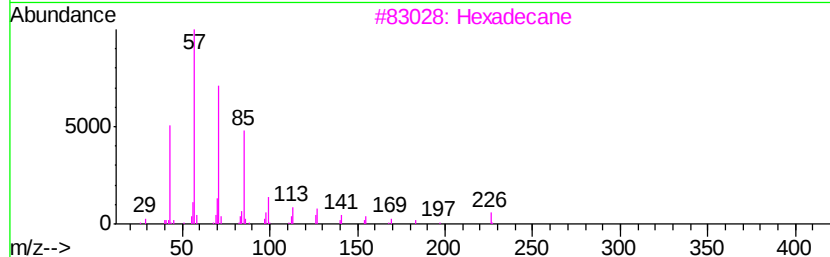
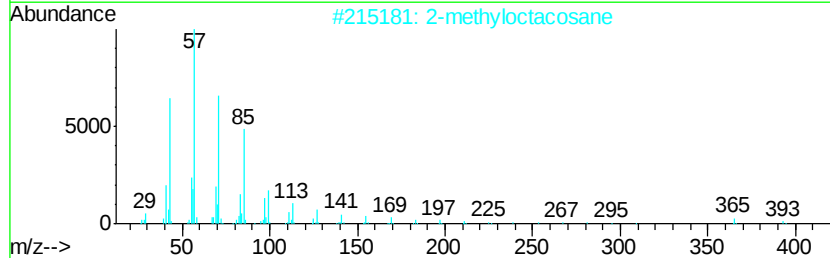
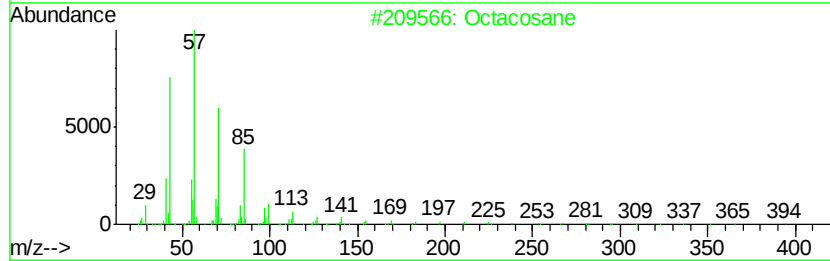
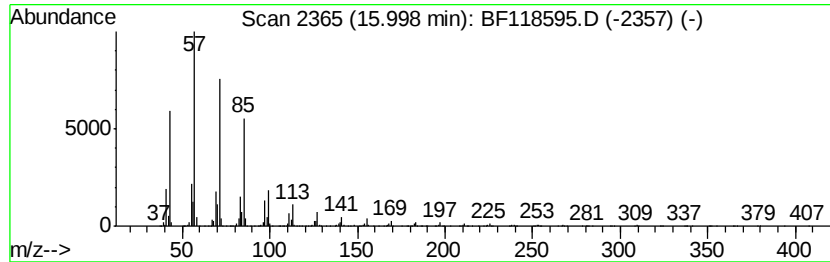
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 2-methyloctacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.00	56.20 ng	8222100	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Hexadecane	226	C16H34	000544-76-3	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118595.D
Acq On : 20 Jan 2020 17:07
Operator : CG/JU
Sample : L1154-07
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B3-15-20

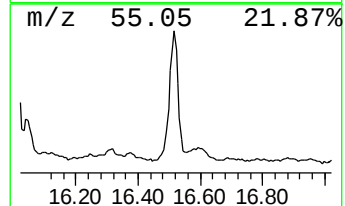
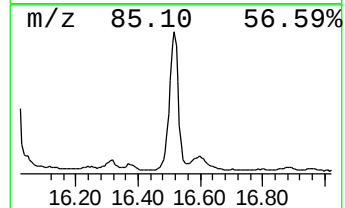
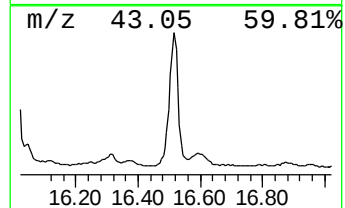
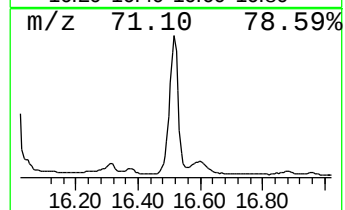
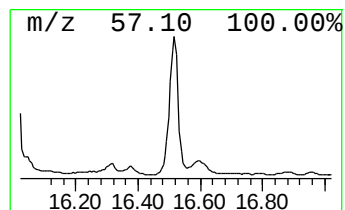
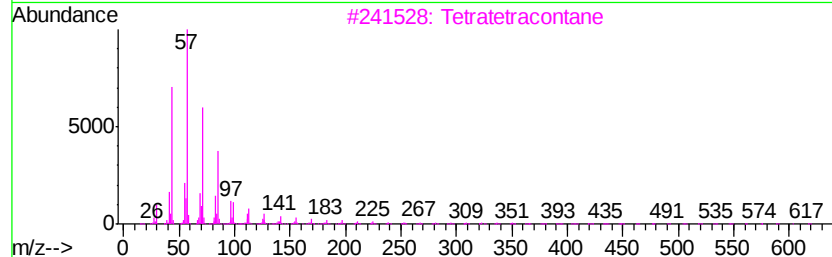
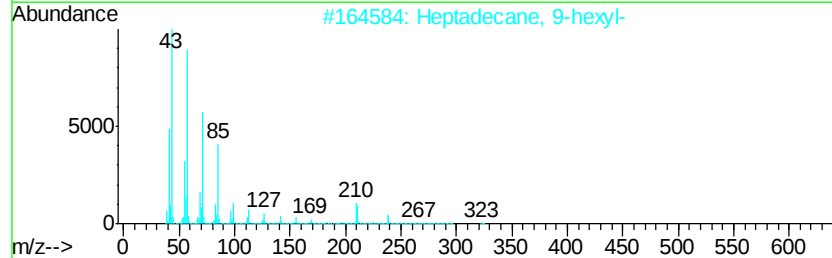
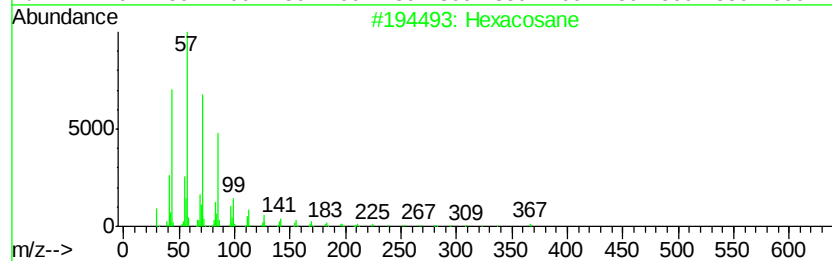
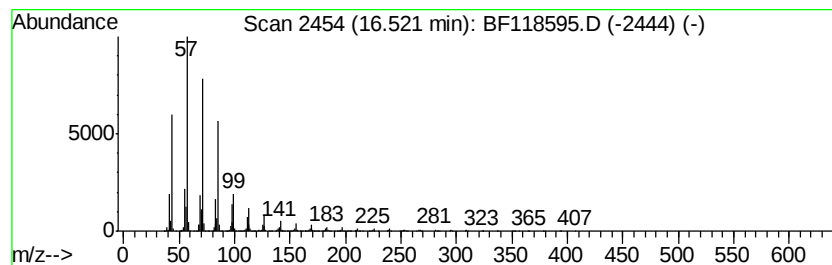
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Heptadecane, 9-hexyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.52	29.25 ng	4279300	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	94
2		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
3		Tetratetracontane	619	C44H90	007098-22-8	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF012020\
 Data File : BF118595.D
 Acq On : 20 Jan 2020 17:07
 Operator : CG/JU
 Sample : L1154-07
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B3-15-20

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012020.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...	2.22	29.5	ng	4037390	1	6.88	2739380	20.0
2-Pentanone, 4-hy...	5.10	10.9	ng	1490470	1	6.88	2739380	20.0
unknown6.65	6.65	78.5	ng	10748900	1	6.88	2739380	20.0
Eicosane, 9-octyl-	12.93	10.9	ng	2285820	5	14.04	4190690	20.0
Eicosane	13.46	266.9	ng	55913200	5	14.04	4190690	20.0
Ethanol, 2-(tetra...	13.89	20.5	ng	4293040	5	14.04	4190690	20.0
Tricosane	13.98	18.1	ng	3785560	5	14.04	4190690	20.0
Tetracosane	14.29	136.3	ng	28555300	5	14.04	4190690	20.0
Heneicosane	14.52	42.1	ng	8815110	5	14.04	4190690	20.0
Hexadecane	14.83	79.4	ng	11612200	6	15.51	2926190	20.0
Nonadecane	14.95	59.3	ng	8672700	6	15.51	2926190	20.0
Heptacosane	15.17	85.3	ng	12486200	6	15.51	2926190	20.0
Octacosane	15.71	14.5	ng	2121990	6	15.51	2926190	20.0
2-methyloctacosane	16.00	56.2	ng	8222100	6	15.51	2926190	20.0
Heptadecane, 9-he...	16.52	29.3	ng	4279300	6	15.51	2926190	20.0