

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
 Data File : BF109909.D
 Acq On : 16 Oct 2018 16:59
 Operator : JU/SJ
 Sample : J5516-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSSI-1011-S-DH-32

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-BF101518.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	2.116	3	5	12	rVB	371221	543398	13.30%	1.277%
2	2.328	34	41	48	rVB	805943	1055975	25.85%	2.481%
3	3.587	254	255	266	rBV4	43207	85396	2.09%	0.201%
4	5.216	523	532	539	rBV	164098	190845	4.67%	0.448%
5	5.598	592	597	601	rBV	4091296	4035748	98.81%	9.482%
6	6.598	761	767	770	rBV	3937671	3948641	96.67%	9.278%
7	6.745	788	792	796	rBV	4301572	4084455	100.00%	9.597%
8	6.981	828	832	835	rVB	1729167	1342235	32.86%	3.154%
9	7.139	855	859	862	rVB	3603556	3318097	81.24%	7.796%
10	7.539	922	927	930	rBV	2931034	2492679	61.03%	5.857%
11	8.263	1046	1050	1053	rBV	2004711	1589459	38.91%	3.735%
12	9.333	1226	1232	1236	rBV	4337278	3915236	95.86%	9.199%
13	9.439	1247	1250	1253	rVB	160828	133436	3.27%	0.314%
14	9.727	1295	1299	1302	rBV	376541	300865	7.37%	0.707%
15	10.022	1345	1349	1353	rBV	1700219	1408024	34.47%	3.308%
16	10.810	1479	1483	1486	rBV	2119720	1761791	43.13%	4.139%
17	11.510	1599	1602	1606	rBV	1283000	1197495	29.32%	2.814%
18	12.016	1685	1688	1694	rBV2	71303	89957	2.20%	0.211%
19	12.727	1806	1809	1814	rBV	65017	68828	1.69%	0.162%
20	12.904	1836	1839	1842	rBV	50443	44850	1.10%	0.105%
21	12.957	1845	1848	1850	rBV	79261	58590	1.43%	0.138%
22	13.098	1868	1872	1875	rBV	3995596	3307974	80.99%	7.772%
23	13.315	1906	1909	1913	rVB	63337	56426	1.38%	0.133%
24	13.568	1947	1952	1956	rBV	384942	527885	12.92%	1.240%
25	13.604	1956	1958	1962	rVV	162202	176228	4.31%	0.414%
26	13.657	1962	1967	1970	rVB2	111977	159531	3.91%	0.375%
27	13.980	2018	2022	2035	rVB2	340855	413386	10.12%	0.971%
28	14.157	2048	2052	2055	rBV	1856275	1659427	40.63%	3.899%
29	14.251	2065	2068	2074	rBV2	104741	166553	4.08%	0.391%
30	14.304	2074	2077	2080	rVB	192897	174160	4.26%	0.409%
31	14.609	2125	2129	2136	rVB	294208	311369	7.62%	0.732%
32	14.927	2180	2183	2190	rVB	287644	318722	7.80%	0.749%
33	15.009	2194	2197	2201	rVB	78453	84393	2.07%	0.198%
34	15.227	2230	2234	2241	rBV2	52305	105804	2.59%	0.249%

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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.286	2241	2244	2257	rVB	307701	398479	9.76%	0.936%
36	15.686	2306	2312	2321	rVB	1516968	1996287	48.88%	4.690%
37	16.156	2387	2392	2397	rBV	236845	357349	8.75%	0.840%
38	16.203	2398	2400	2409	rVB9	40511	81059	1.98%	0.190%
39	16.698	2479	2484	2491	rVB	162121	288691	7.07%	0.678%
40	17.339	2589	2593	2600	rVB2	75037	142744	3.49%	0.335%
41	17.868	2679	2683	2691	rVB10	78060	168545	4.13%	0.396%

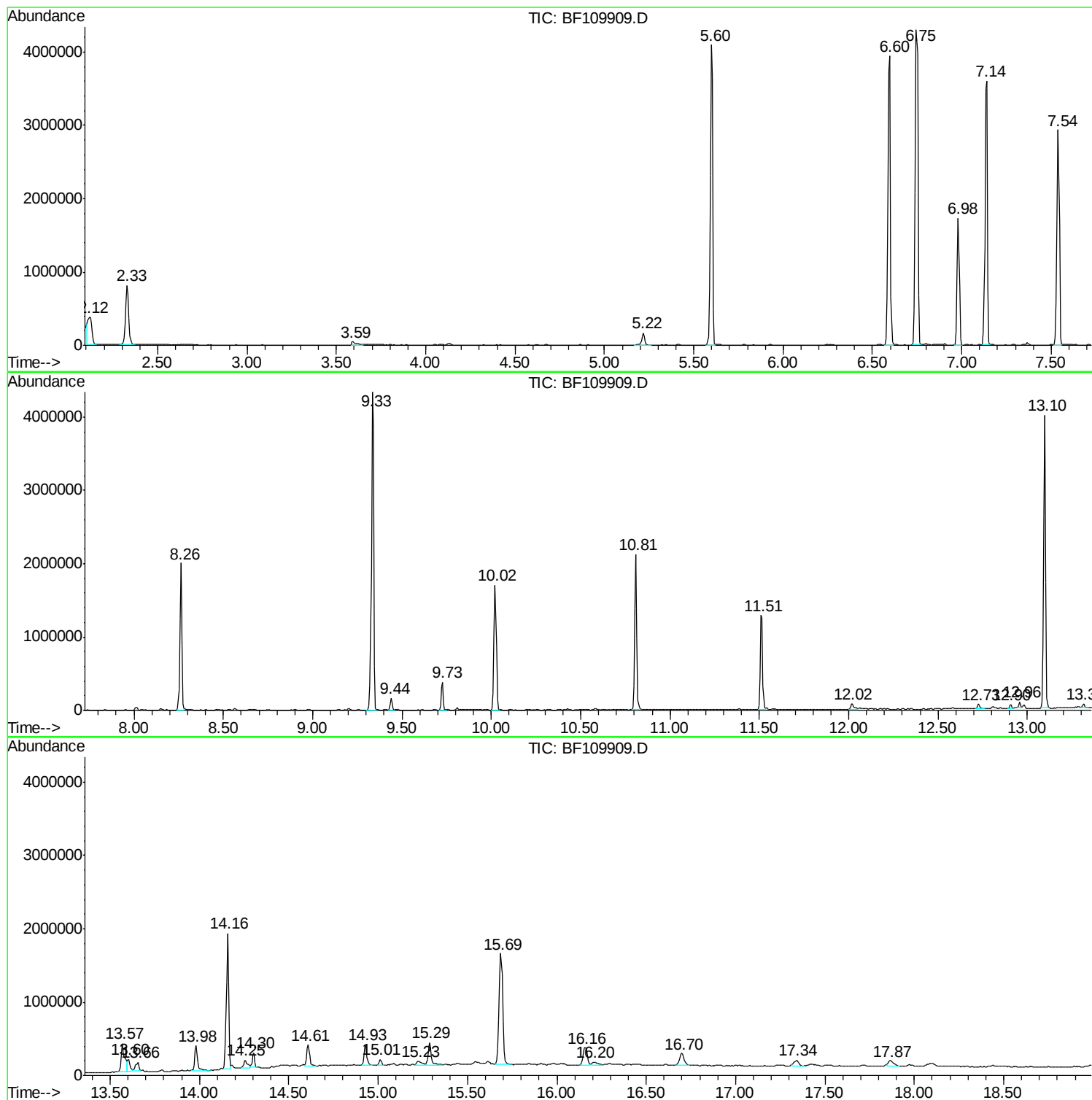
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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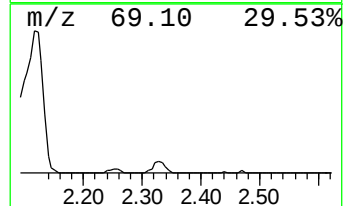
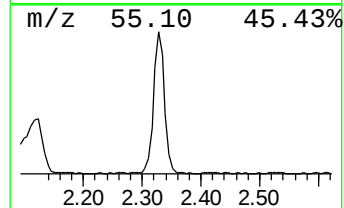
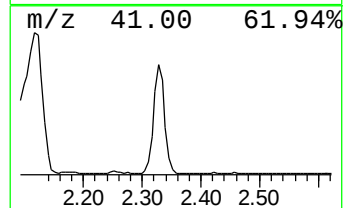
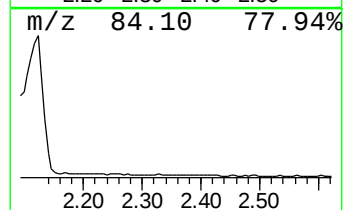
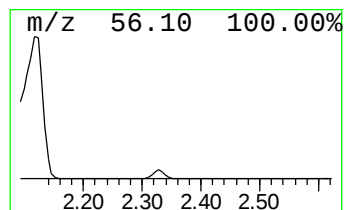
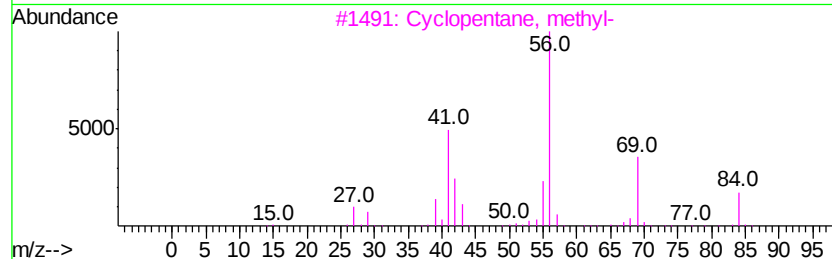
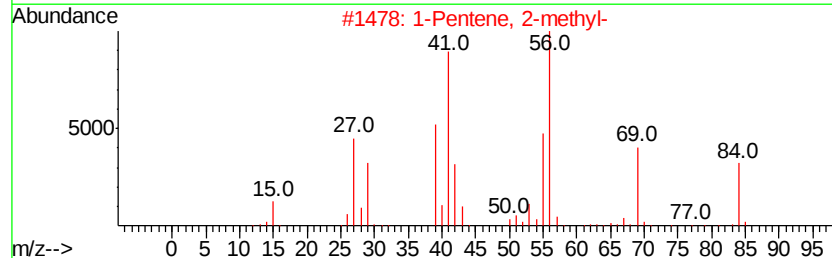
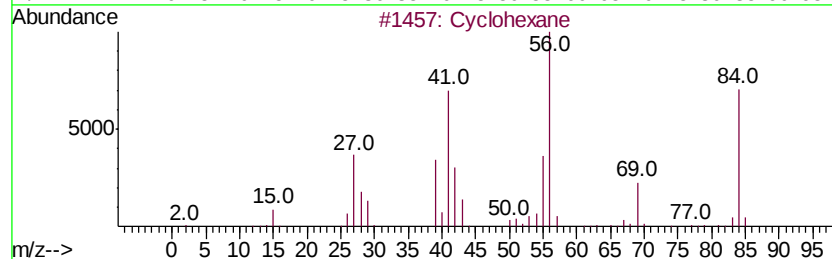
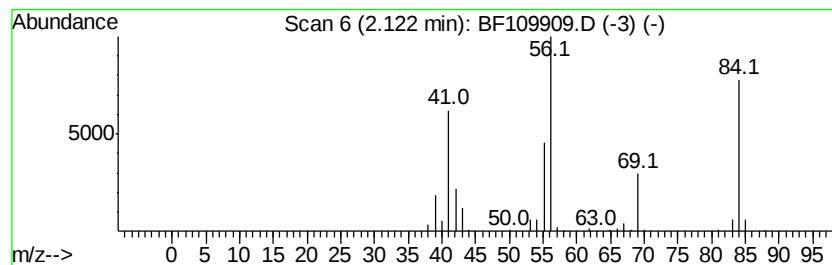
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclohexane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.12	8.10 ng	543398	1,4-Dichlorobenzene-d4	6.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane	84	C6H12	000110-82-7	91
2		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	78
3		Cyclopentane, methyl-	84	C6H12	000096-37-7	56
4		Cyclopropane, propyl-	84	C6H12	002415-72-7	53
5		Cyclobutane	56	C4H8	000287-23-0	43



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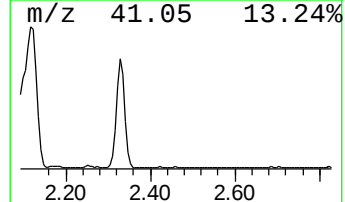
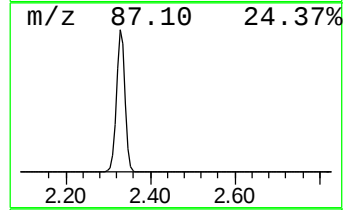
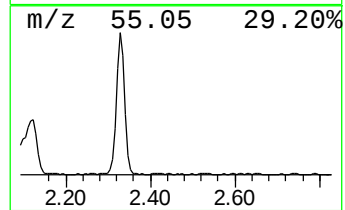
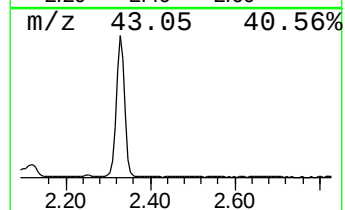
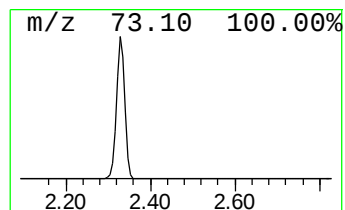
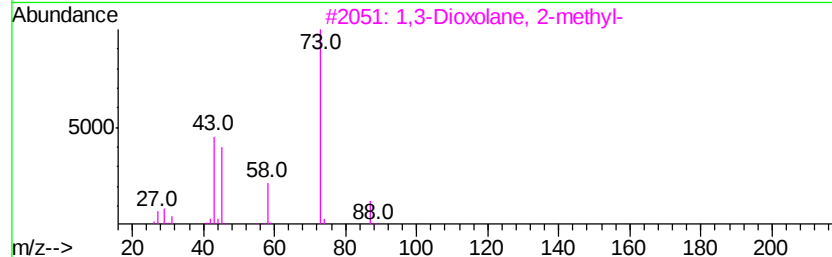
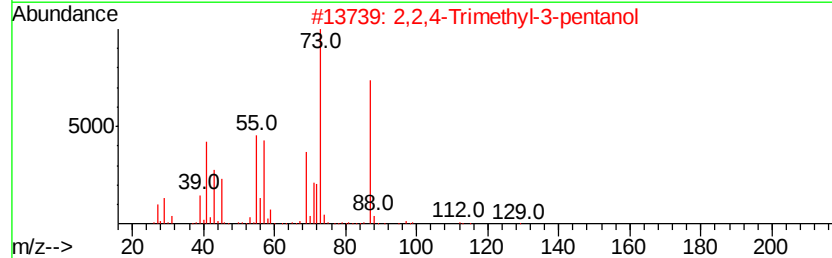
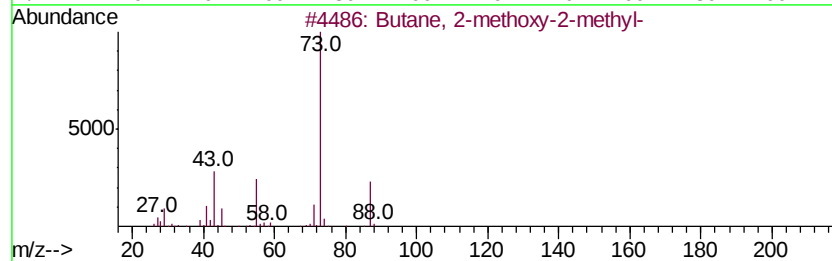
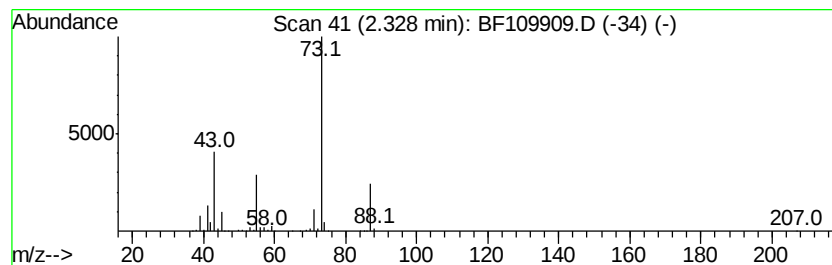
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Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.33	15.73 ng	1055980	1,4-Dichlorobenzene-d4	6.98

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	10



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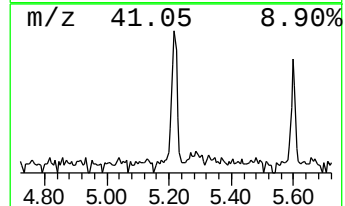
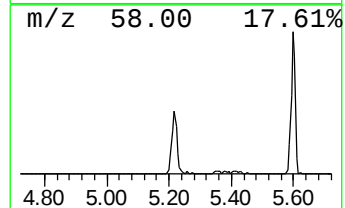
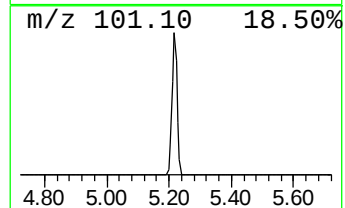
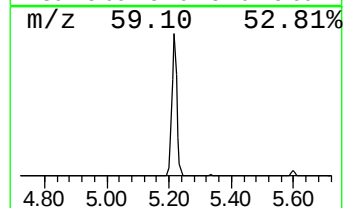
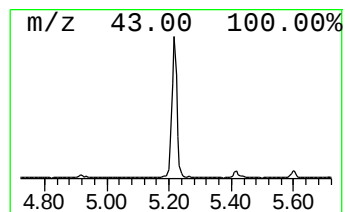
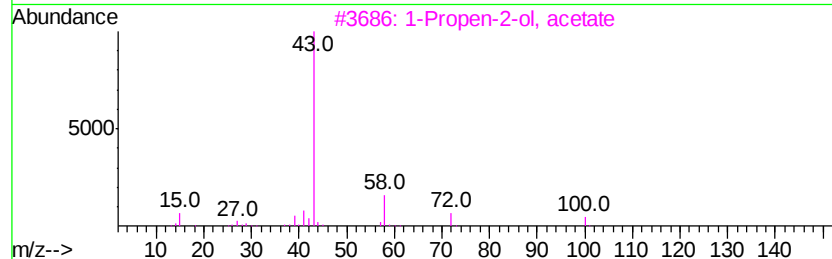
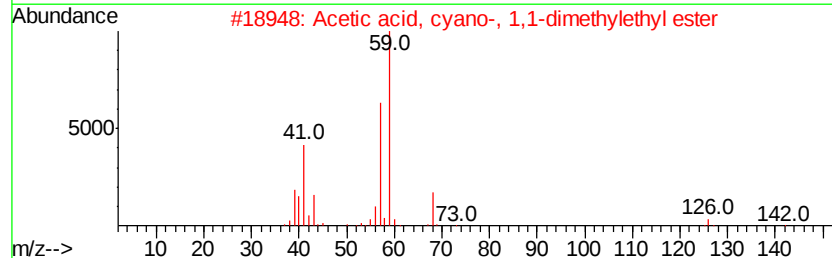
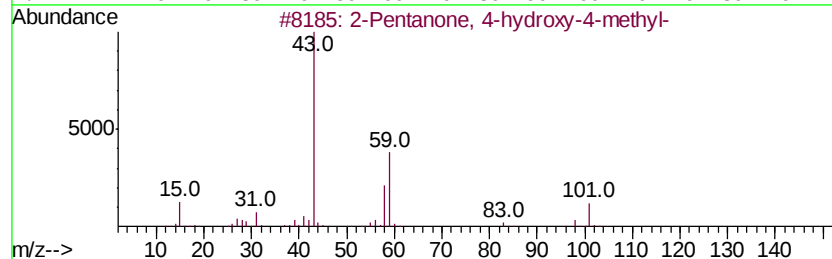
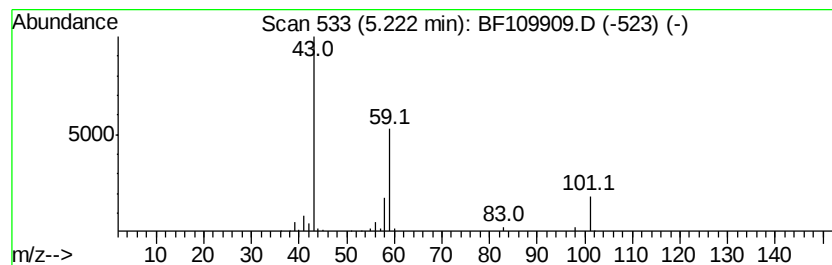
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	2.84 ng	190845	1,4-Dichlorobenzene-d4	6.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4			1,2-Butanediol	90	C4H10O2	000584-03-2	9
5			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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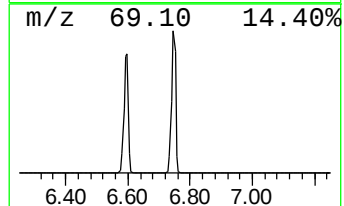
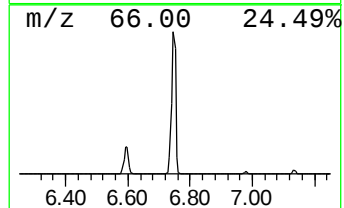
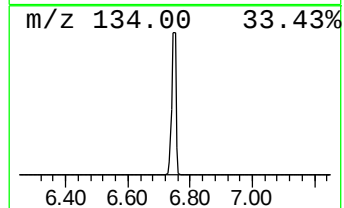
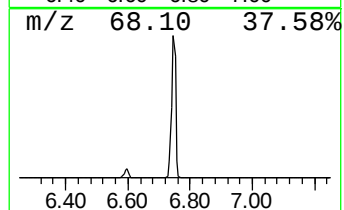
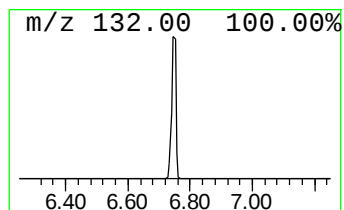
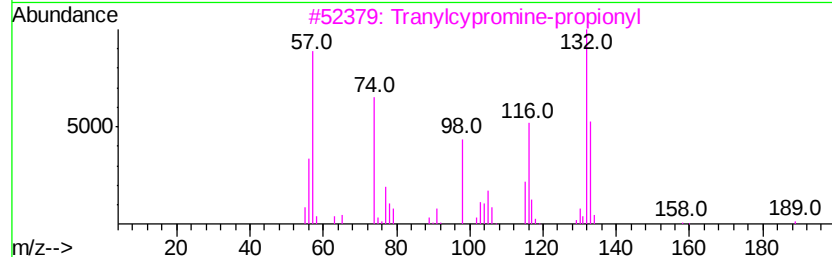
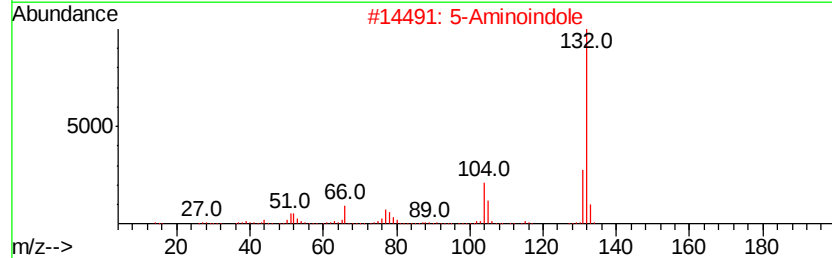
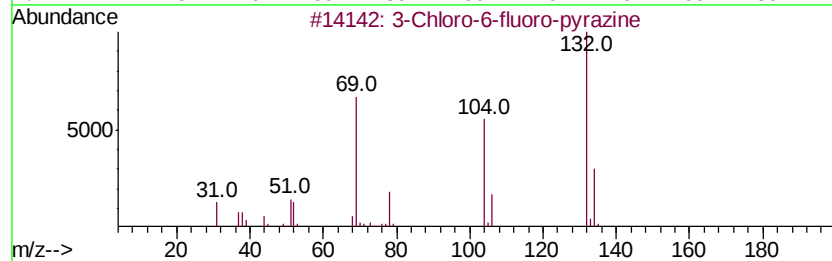
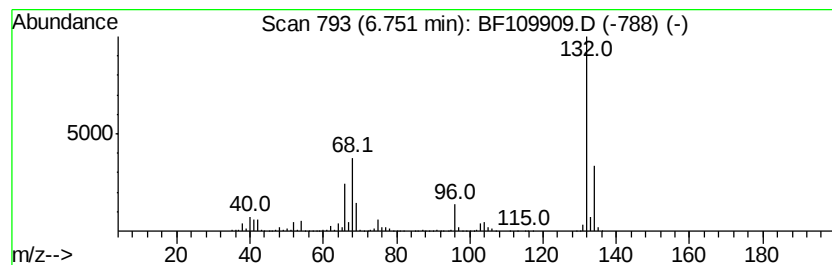
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.75 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	60.86 ng	4084460	1,4-Dichlorobenzene-d4	6.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27	
2	5-Aminoindole	132	C8H8N2	005192-03-0	12	
3	Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10	
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9	
5	4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9	



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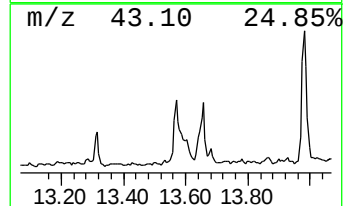
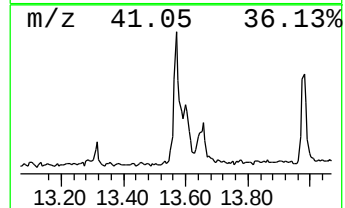
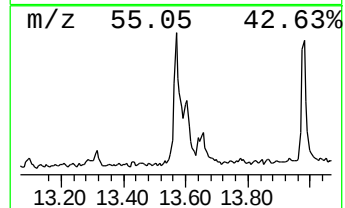
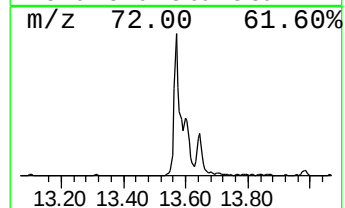
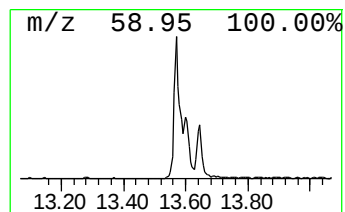
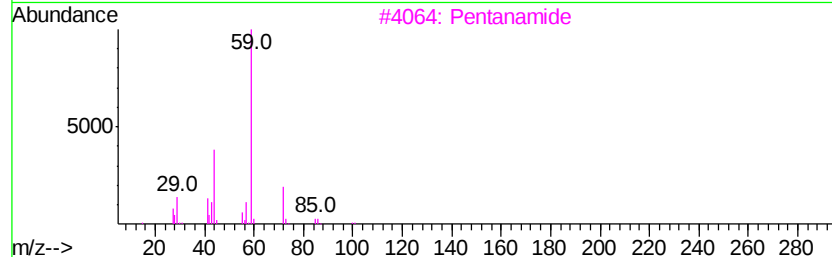
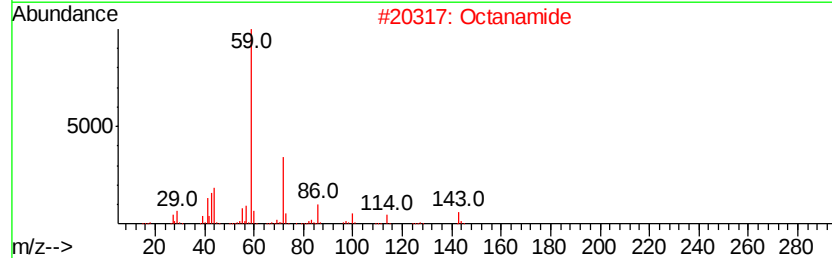
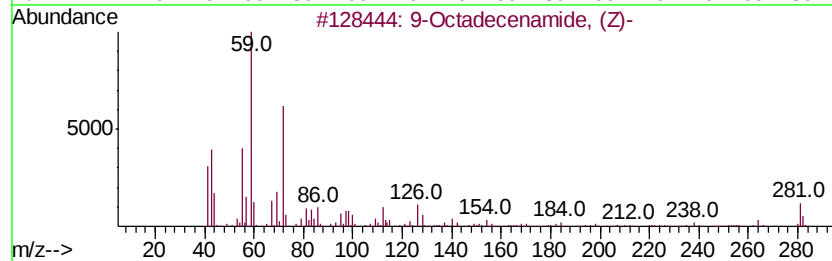
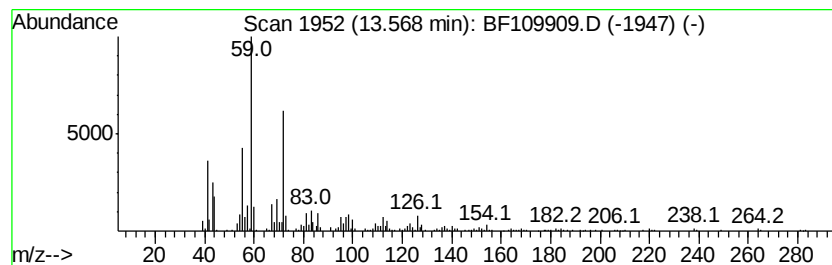
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 9-Octadecenamide, (Z)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.57	6.36 ng	527885	Chrysene-d12	14.16

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	94
2		Octanamide	143	C8H17NO	000629-01-6	59
3		Pentanamide	101	C5H11NO	000626-97-1	59
4		Nonanamide	157	C9H19NO	001120-07-6	59
5		Hexadecanamide	255	C16H33NO	000629-54-9	56



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
Data File : BF109909.D
Acq On : 16 Oct 2018 16:59
Operator : JU/SJ
Sample : J5516-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SSSI-1011-S-DH-32

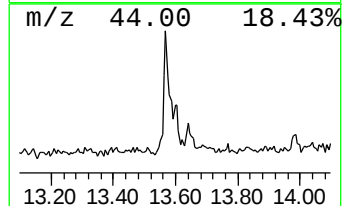
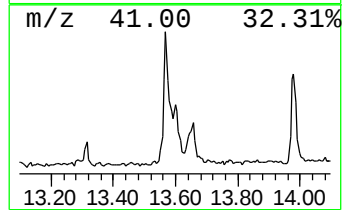
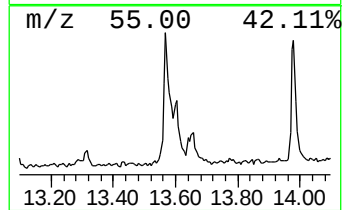
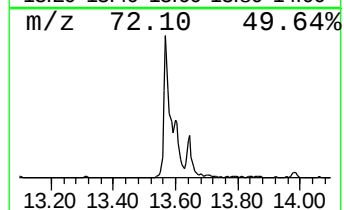
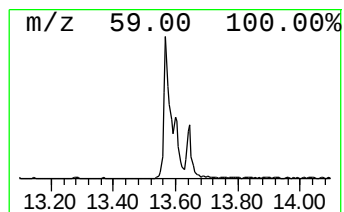
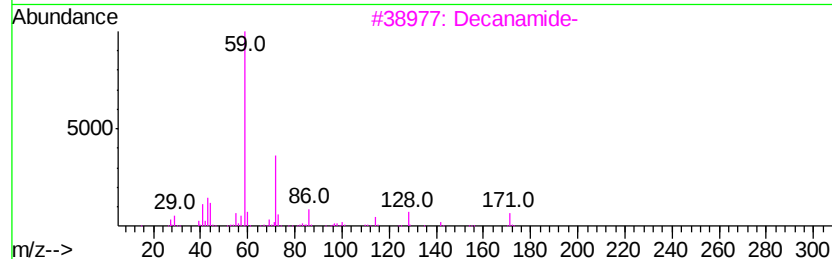
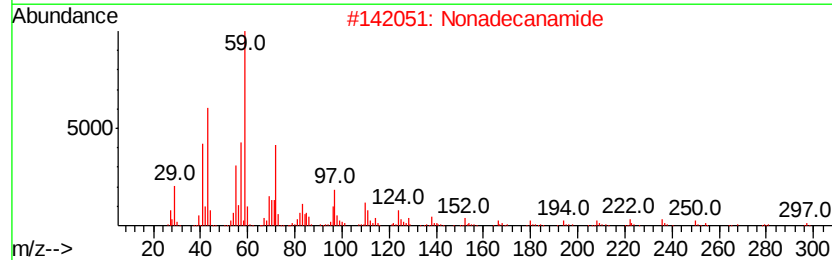
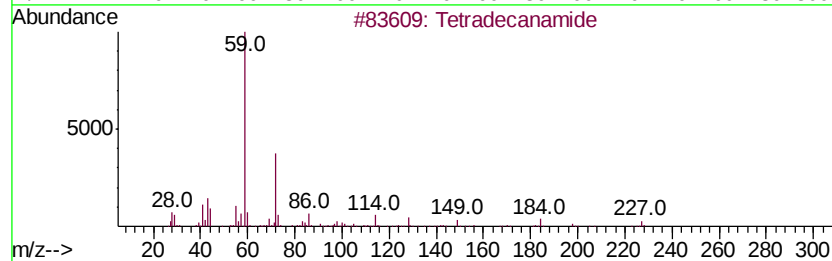
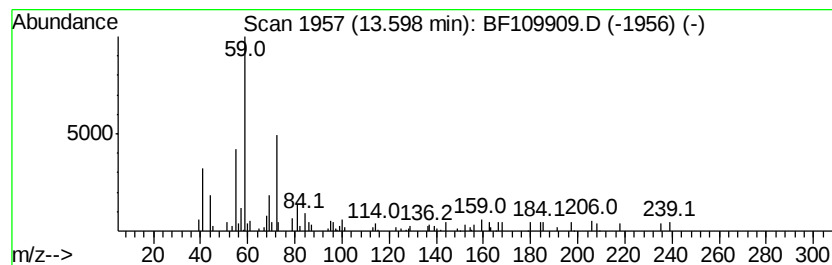
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetradecanamide Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.60	2.12 ng	176228	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecanamide	227	C14H29NO	000638-58-4	64
2		Nonadecanamide	297	C19H39NO	058185-32-3	59
3		Decanamide-	171	C10H21NO	002319-29-1	58
4		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	50
5		Octadecanamide	283	C18H37NO	000124-26-5	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
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ClientSampleId :
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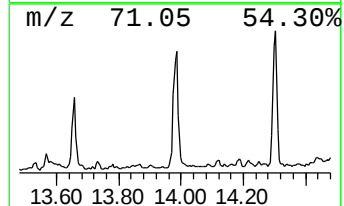
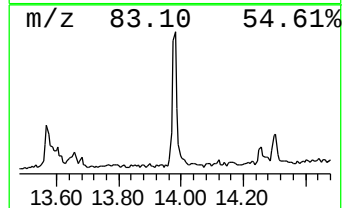
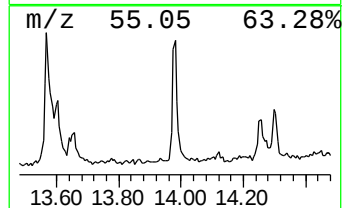
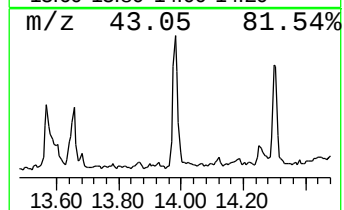
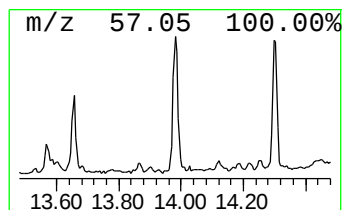
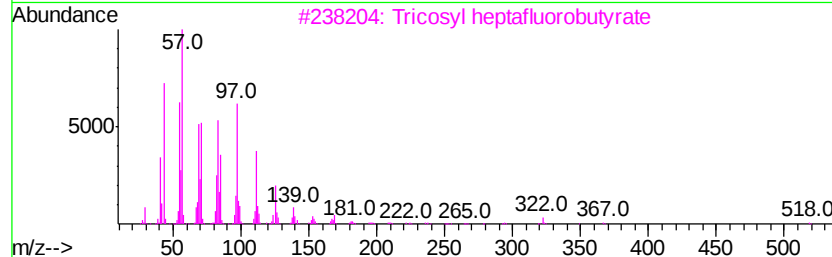
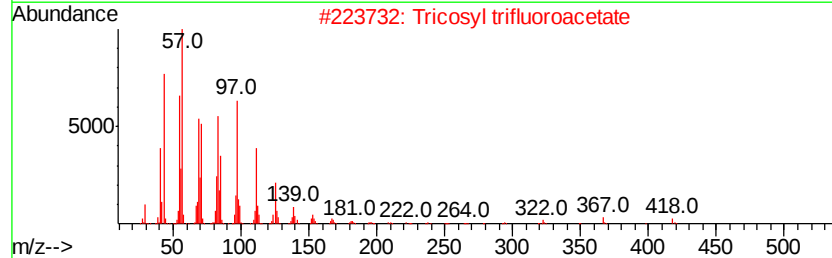
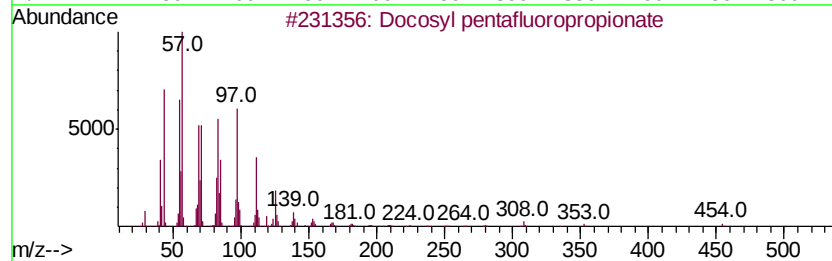
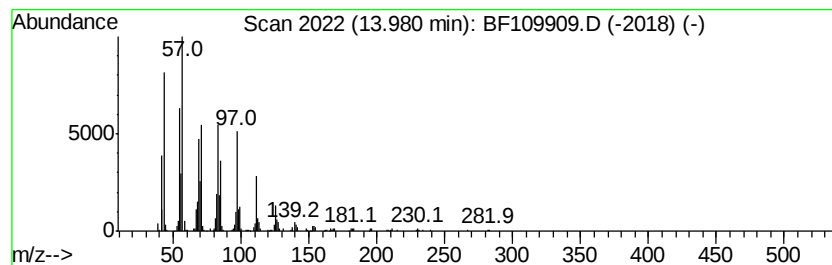
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Docosyl pentafluoropropionate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.98	4.98 ng	413386	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	91
2		Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	91
3		Tricosyl heptafluorobutrate	536	C27H47F7O2	1000351-83-4	91
4		Tetracosyl heptafluorobutrate	550	C28H49F7O2	1000351-83-7	91
5		Tricosyl pentafluoropropionate	486	C26H47F5O2	1000351-81-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
Data File : BF109909.D
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Operator : JU/SJ
Sample : J5516-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

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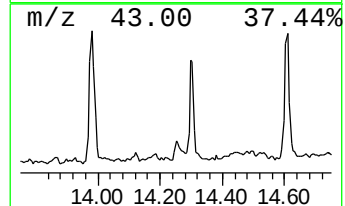
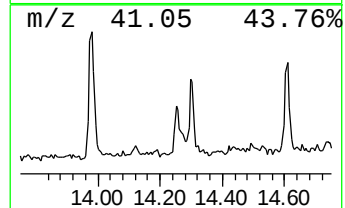
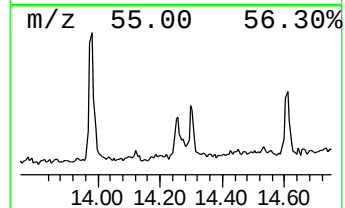
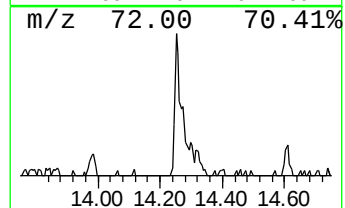
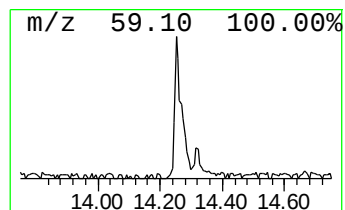
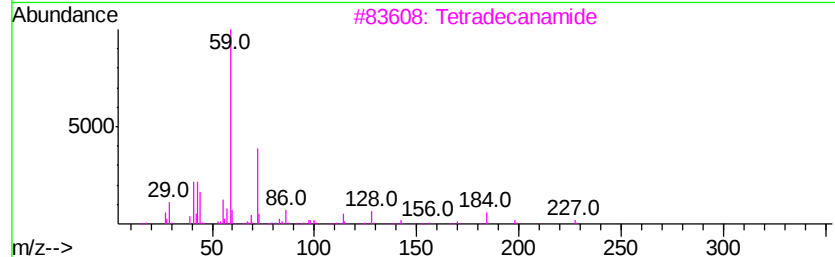
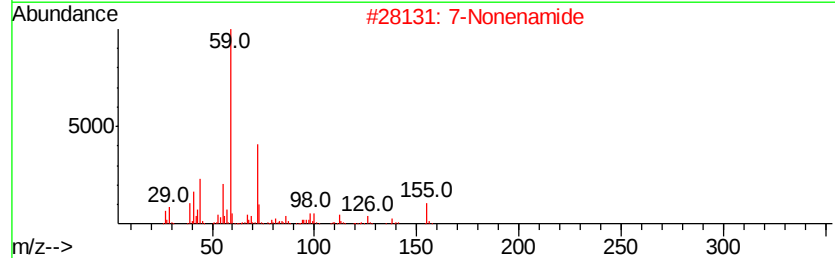
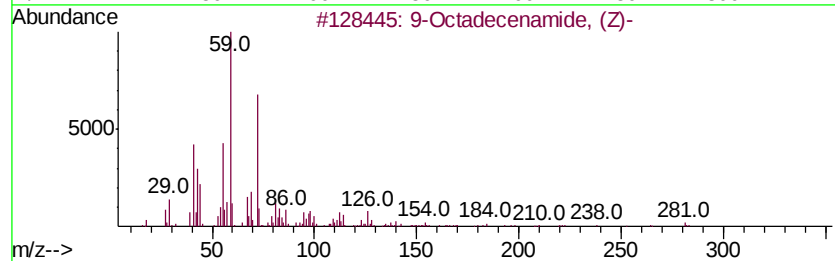
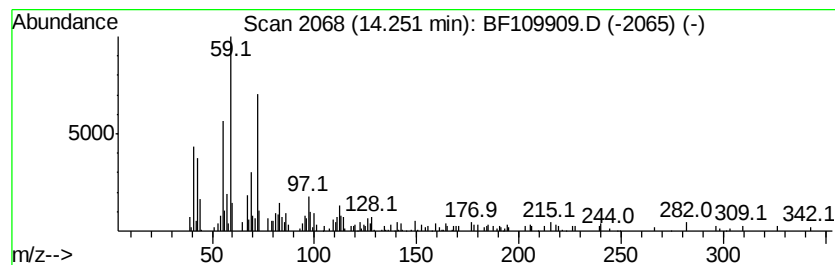
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 7-Nonenamide Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.25	2.01 ng	166553	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	87
2		7-Nonenamide	155	C9H17NO	090949-53-4	72
3		Tetradecanamide	227	C14H29NO	000638-58-4	68
4		Pentanamide, 4-methyl-	115	C6H13NO	001119-29-5	50
5		Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
Data File : BF109909.D
Acq On : 16 Oct 2018 16:59
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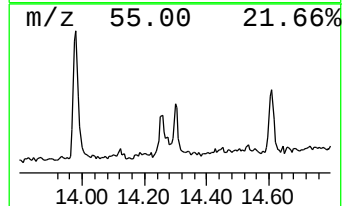
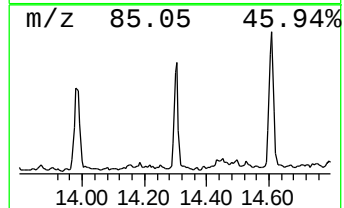
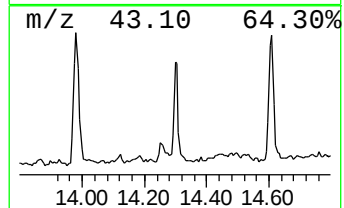
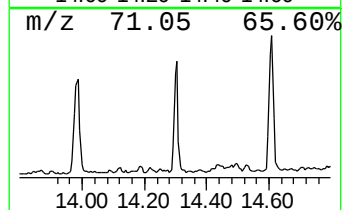
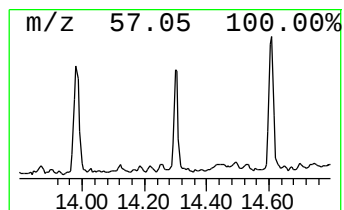
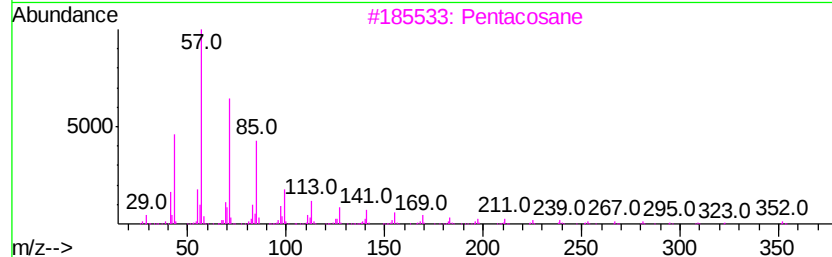
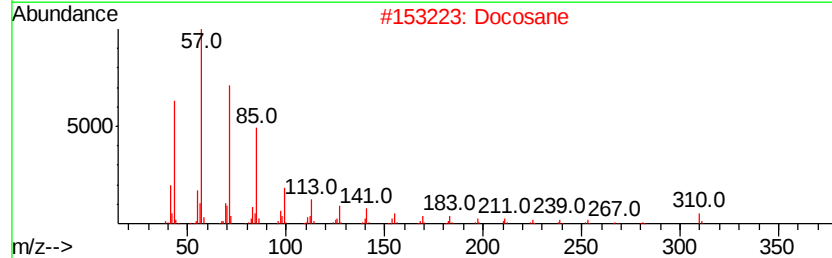
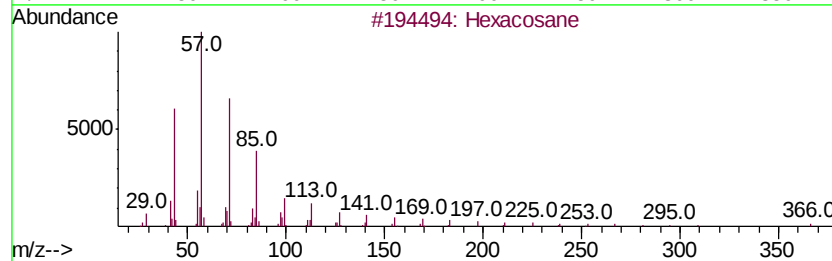
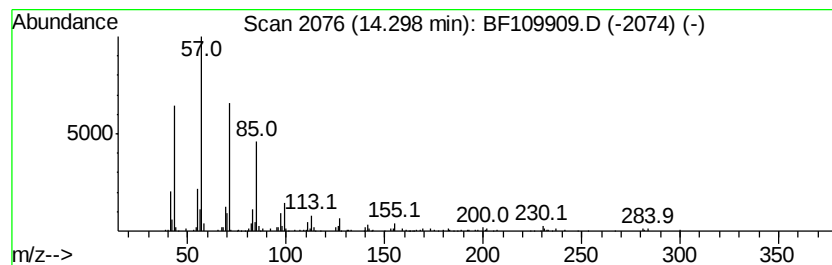
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.30	2.10 ng	174160	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Docosane	310	C22H46	000629-97-0	90
3		Pentacosane	352	C25H52	000629-99-2	83
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	80
5		Octadecane	254	C18H38	000593-45-3	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
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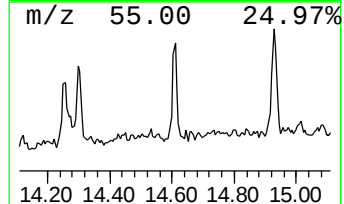
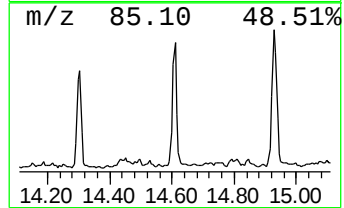
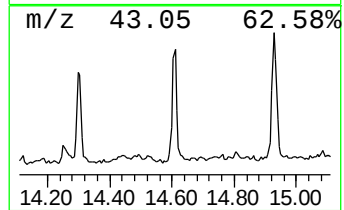
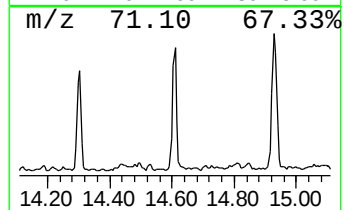
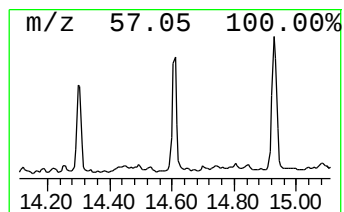
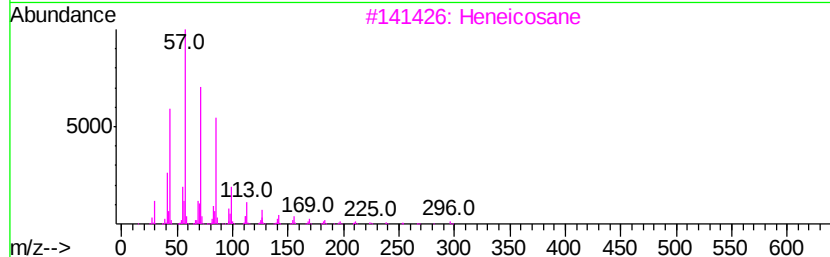
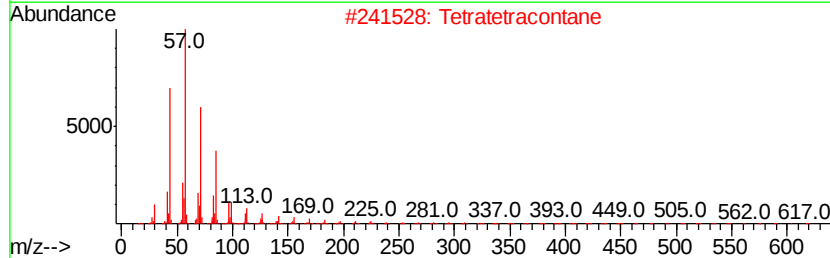
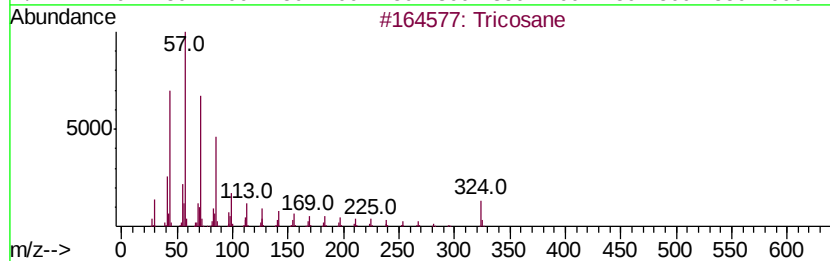
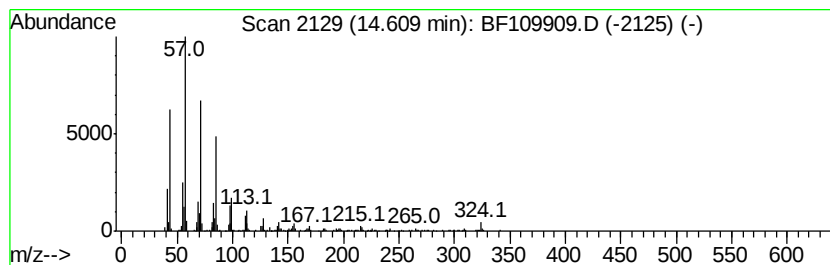
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tricosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.61	3.75 ng	311369	Chrysene-d12	14.16	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricosane	324	C23H48	000638-67-5	91
2	Tetratetracontane	619	C44H90	007098-22-8	90
3	Heneicosane	296	C21H44	000629-94-7	90
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	87



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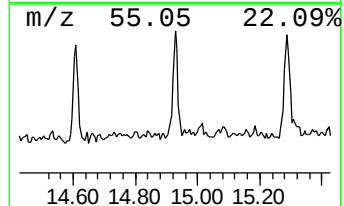
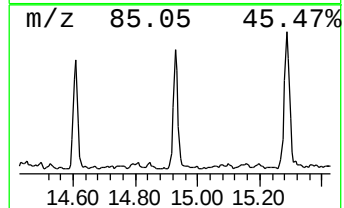
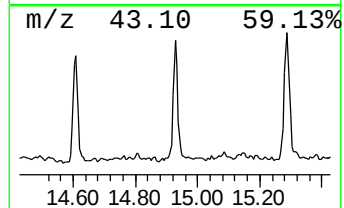
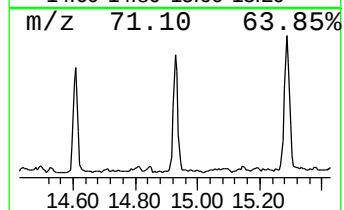
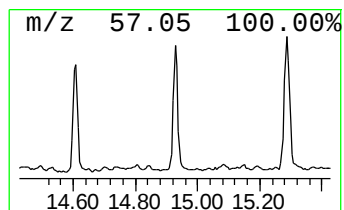
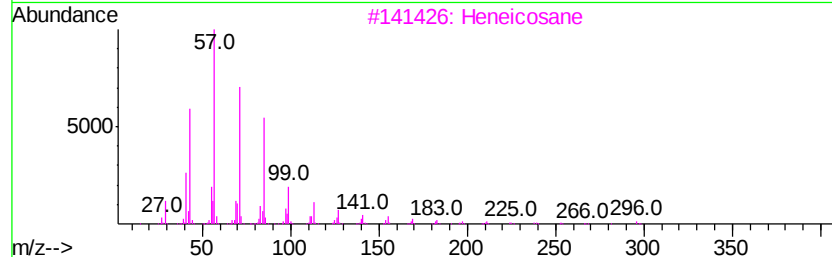
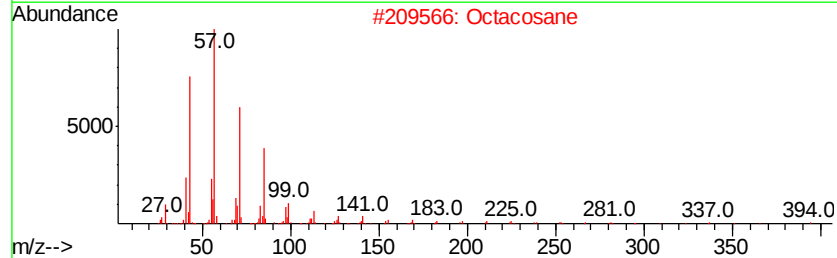
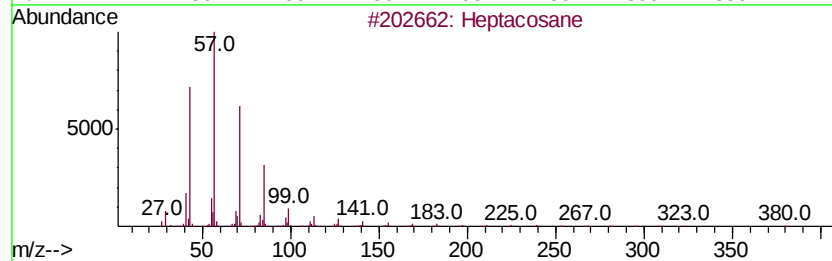
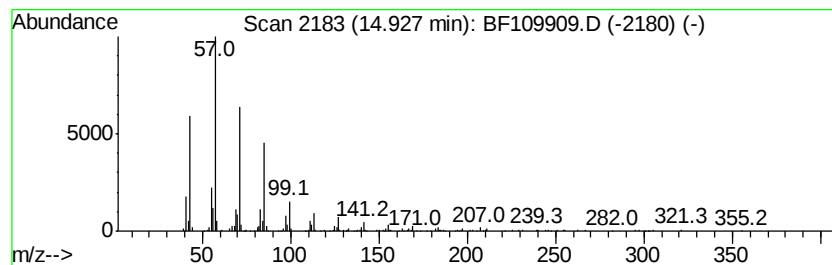
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.93	3.19 ng	318722	Perylene-d12	15.69
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptacosane	380 C27H56	000593-49-7	91
2	Octacosane	394 C28H58	000630-02-4	91
3	Heneicosane	296 C21H44	000629-94-7	91
4	Docosane	310 C22H46	000629-97-0	91
5	Docosane, 7-hexyl-	394 C28H58	055373-86-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
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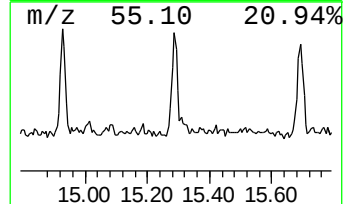
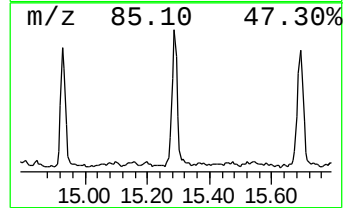
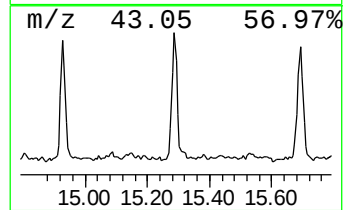
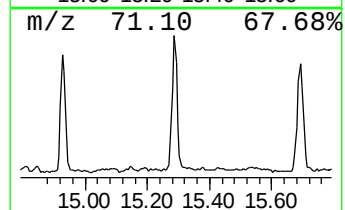
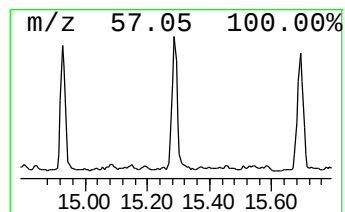
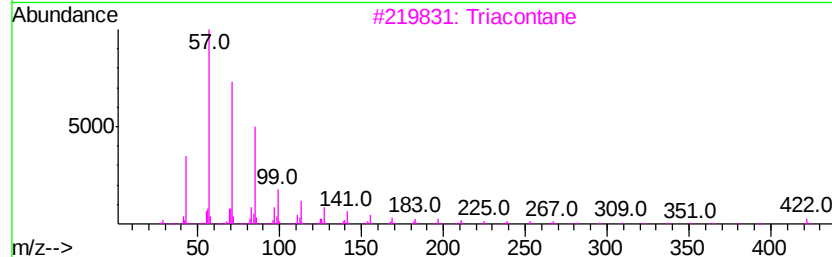
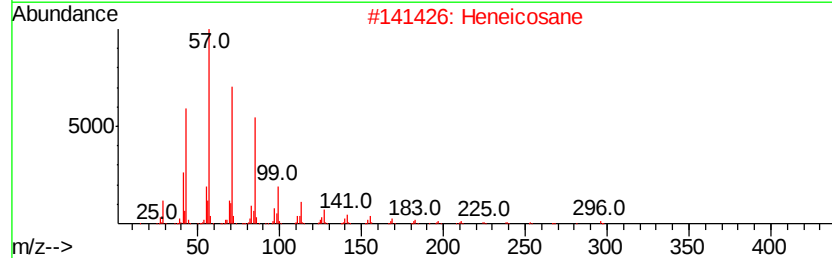
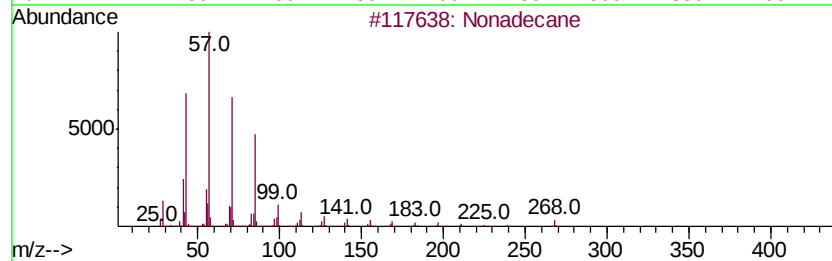
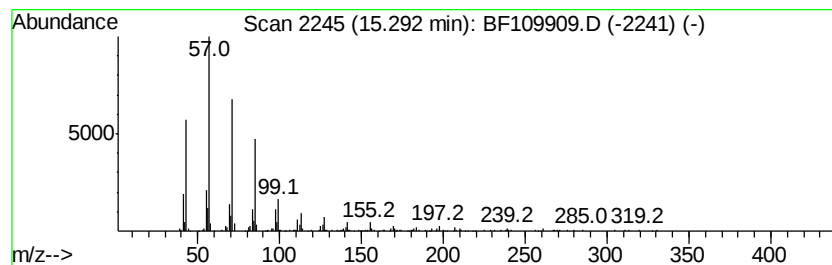
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Nonadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.29	3.99 ng	398479	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	93
2		Heneicosane	296	C21H44	000629-94-7	91
3		Triacontane	422	C30H62	000638-68-6	91
4		Docosane, 11-butyl-	366	C26H54	013475-76-8	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
Data File : BF109909.D
Acq On : 16 Oct 2018 16:59
Operator : JU/SJ
Sample : J5516-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSSI-1011-S-DH-32

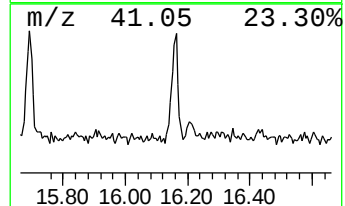
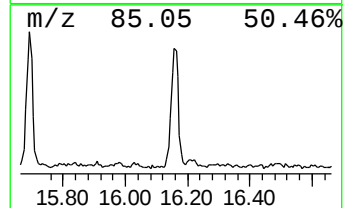
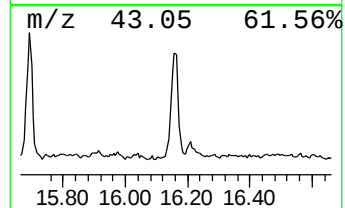
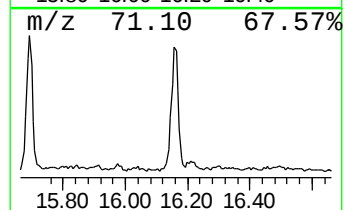
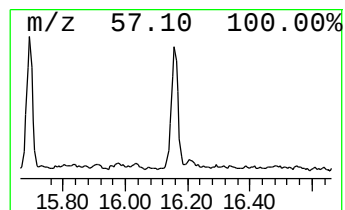
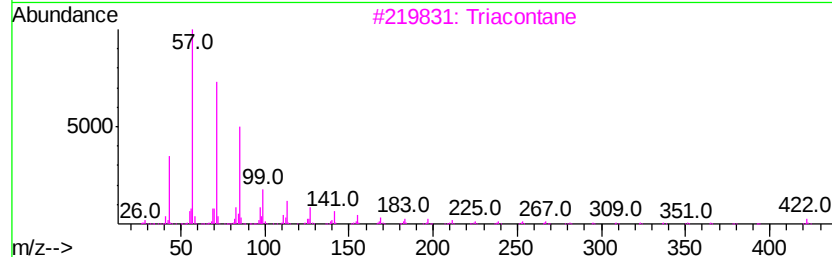
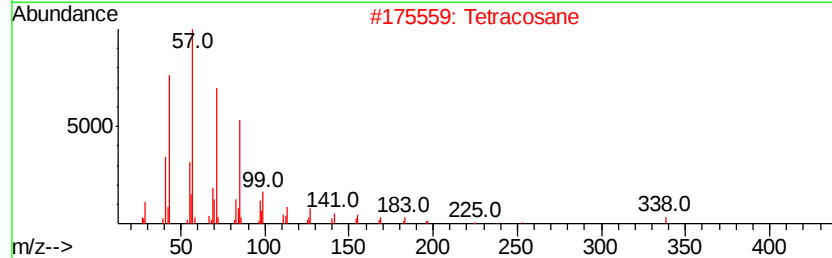
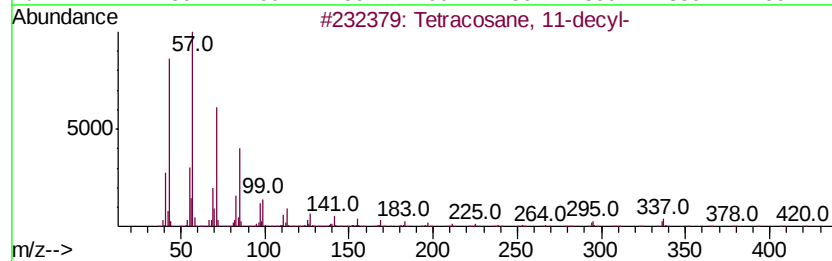
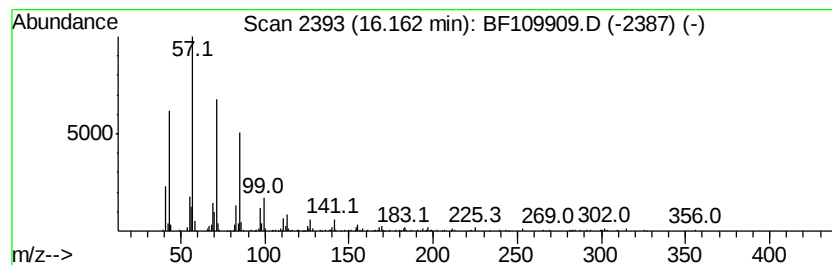
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Tetracosane, 11-decyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	3.58 ng	357349	Perylene-d12	15.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91
2			Tetracosane	338	C24H50	000646-31-1	87
3			Triacontane	422	C30H62	000638-68-6	87
4			Eicosane	282	C20H42	000112-95-8	87
5			2-methylhexacosane	380	C27H56	1000376-72-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
Data File : BF109909.D
Acq On : 16 Oct 2018 16:59
Operator : JU/SJ
Sample : J5516-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSSI-1011-S-DH-32

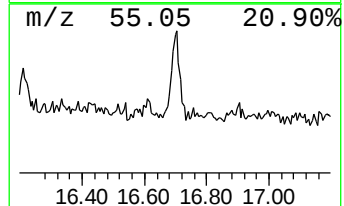
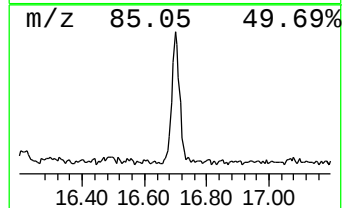
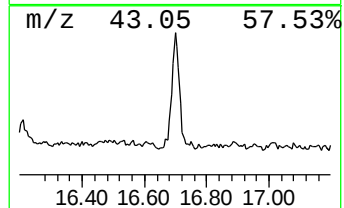
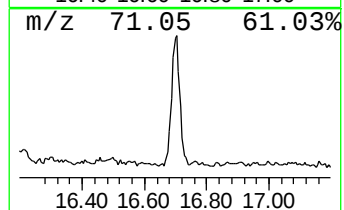
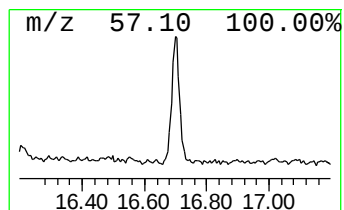
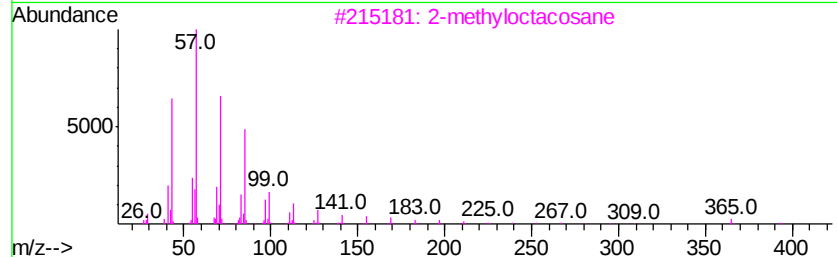
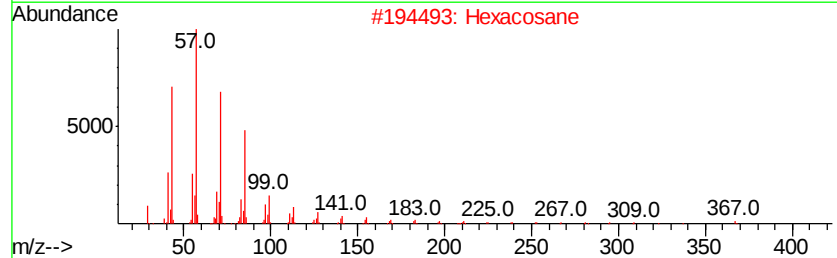
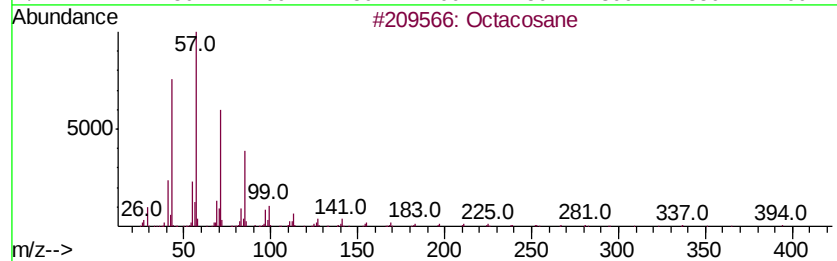
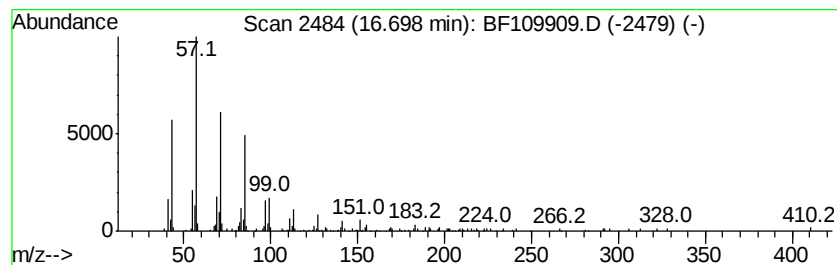
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Octacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.70	2.89 ng	288691	Perylene-d12	15.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	90
2			Hexacosane	366	C26H54	000630-01-3	90
3			2-methyloctacosane	408	C29H60	1000376-72-8	90
4			Tetratetracontane	619	C44H90	007098-22-8	87
5			Hentriacontane	437	C31H64	000630-04-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101618\
Data File : BF109909.D
Acq On : 16 Oct 2018 16:59
Operator : JU/SJ
Sample : J5516-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSSI-1011-S-DH-32

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101518.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
Cyclohexane	2.12	8.1 ng		543398	1	6.98	1342240	20.0
Butane, 2-methoxy...	2.33	15.7 ng		1055980	1	6.98	1342240	20.0
2-Pentanone, 4-hy...	5.22	2.8 ng		190845	1	6.98	1342240	20.0
unknown6.75	6.75	60.9 ng		4084460	1	6.98	1342240	20.0
9-Octadecenamide,...	13.57	6.4 ng		527885	5	14.16	1659430	20.0
Tetradecanamide	13.60	2.1 ng		176228	5	14.16	1659430	20.0
Docosyl pentaflu...	13.98	5.0 ng		413386	5	14.16	1659430	20.0
7-Nonenamide	14.25	2.0 ng		166553	5	14.16	1659430	20.0
Hexacosane	14.30	2.1 ng		174160	5	14.16	1659430	20.0
Tricosane	14.61	3.8 ng		311369	5	14.16	1659430	20.0
Heptacosane	14.93	3.2 ng		318722	6	15.69	1996290	20.0
Nonadecane	15.29	4.0 ng		398479	6	15.69	1996290	20.0
Tetracosane, 11-d...	16.16	3.6 ng		357349	6	15.69	1996290	20.0
Octacosane	16.70	2.9 ng		288691	6	15.69	1996290	20.0