

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF120919\
 Data File : BF118149.D
 Acq On : 9 Dec 2019 17:29
 Operator : JU
 Sample : K6187-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RPC-BOT-S

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-BF120619.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	5.075	503	508	519	rVB	372318	445052	17.43%	1.896%
2	5.481	573	577	590	rBV	2075740	1995697	78.14%	8.501%
3	6.498	745	750	753	rBV	2297564	2069046	81.02%	8.814%
4	6.639	769	774	777	rBV	2587209	2286357	89.53%	9.740%
5	6.869	809	813	816	rBV	729024	597144	23.38%	2.544%
6	7.022	835	839	843	rBV	2069266	1810337	70.89%	7.712%
7	7.428	904	908	920	rBV	1387578	1340885	52.50%	5.712%
8	8.151	1028	1031	1050	rVB	844011	768997	30.11%	3.276%
9	9.233	1211	1215	1218	rBV	2863145	2553843	100.00%	10.879%
10	9.627	1279	1282	1290	rVB	191254	189931	7.44%	0.809%
11	9.916	1327	1331	1343	rVB	982097	871942	34.14%	3.714%
12	10.704	1461	1465	1477	rBV	1913145	1723133	67.47%	7.340%
13	11.404	1581	1584	1587	rBV	859762	784231	30.71%	3.341%
14	11.427	1587	1588	1592	rVB	55062	36608	1.43%	0.156%
15	11.898	1665	1668	1670	rBV2	27577	36440	1.43%	0.155%
16	11.927	1670	1673	1687	rVB	336619	363831	14.25%	1.550%
17	12.621	1788	1791	1795	rBV	85259	85761	3.36%	0.365%
18	12.715	1804	1807	1821	rBV3	82583	101968	3.99%	0.434%
19	12.857	1825	1831	1834	rVB	58194	66496	2.60%	0.283%
20	12.998	1850	1855	1858	rBV	2416102	2132782	83.51%	9.085%
21	13.563	1947	1951	1955	rBV2	24609	29040	1.14%	0.124%
22	13.892	2003	2007	2017	rVB	234981	308029	12.06%	1.312%
23	14.057	2027	2035	2038	rBV	673225	660946	25.88%	2.816%
24	14.210	2058	2061	2067	rVB	60266	63256	2.48%	0.269%
25	14.521	2110	2114	2120	rBV3	105699	137940	5.40%	0.588%
26	14.827	2163	2166	2171	rVB	108425	114831	4.50%	0.489%
27	14.980	2189	2192	2196	rVB2	36044	35435	1.39%	0.151%
28	15.109	2210	2214	2217	rBV	45248	64255	2.52%	0.274%
29	15.174	2221	2225	2228	rBV	152934	174741	6.84%	0.744%
30	15.204	2228	2230	2236	rVB5	34641	58816	2.30%	0.251%
31	15.415	2262	2266	2269	rBV	31439	39159	1.53%	0.167%
32	15.480	2273	2277	2283	rBV	35537	50734	1.99%	0.216%
33	15.551	2283	2289	2297	rBV	578507	908724	35.58%	3.871%
34	16.009	2362	2367	2372	rBV	140749	210780	8.25%	0.898%

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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	16.068	2372	2377	2385	rVV3	43322	84989	3.33%	0.362%
36	16.527	2450	2455	2461	rBV2	64699	114670	4.49%	0.488%
37	17.051	2539	2544	2552	rBV4	22464	51914	2.03%	0.221%
38	17.139	2552	2559	2566	rBV5	32106	74931	2.93%	0.319%
39	17.503	2618	2621	2628	rVB3	16898	31205	1.22%	0.133%

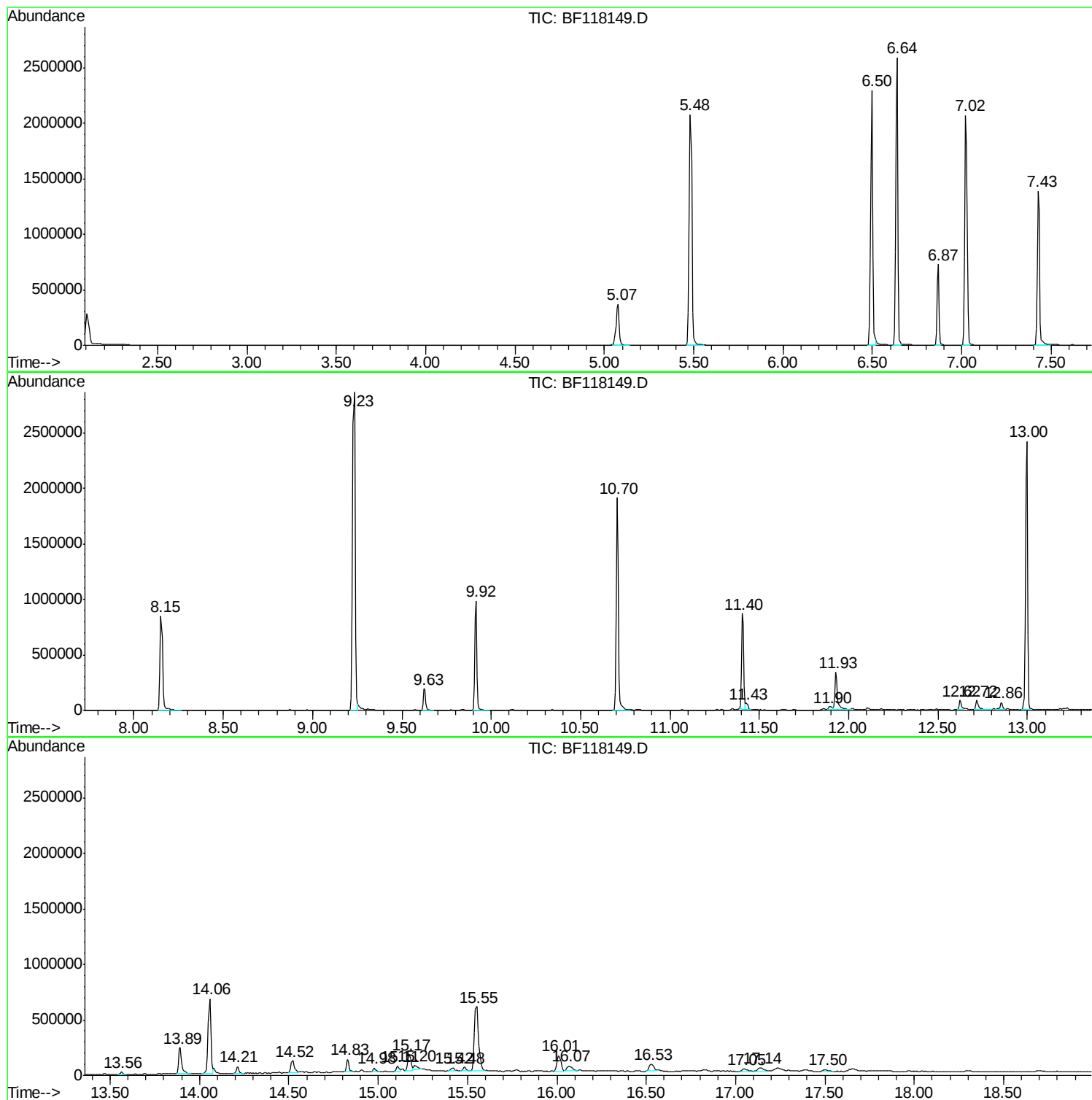
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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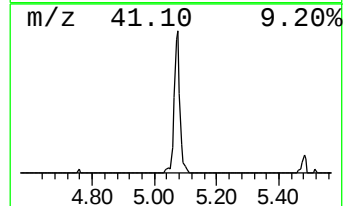
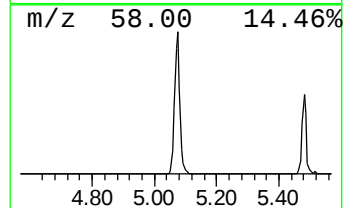
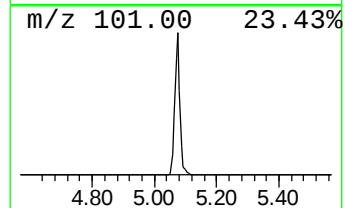
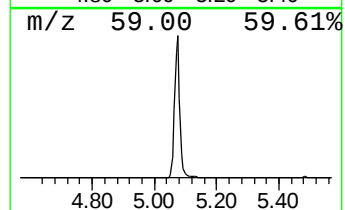
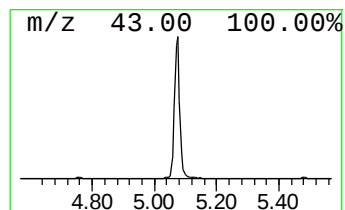
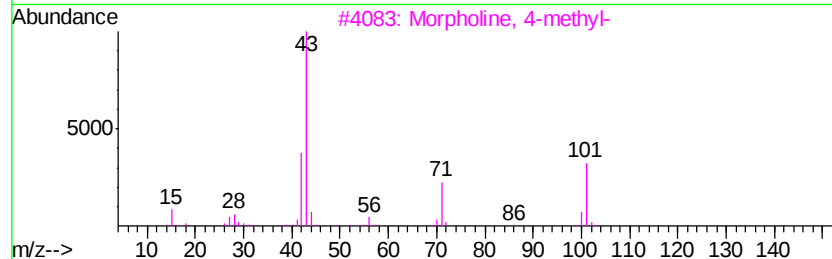
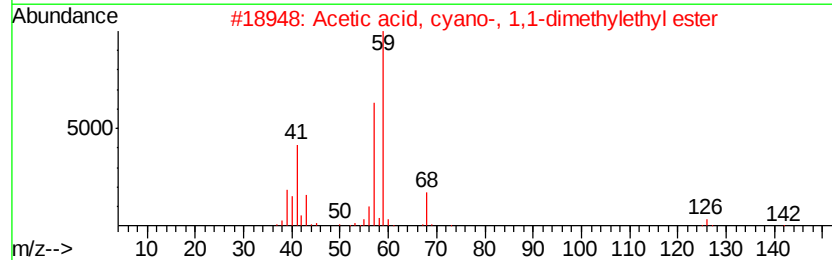
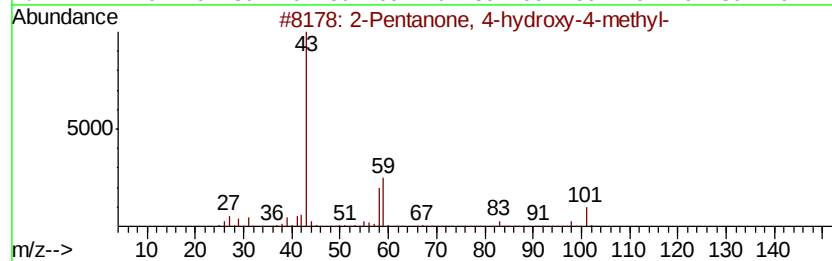
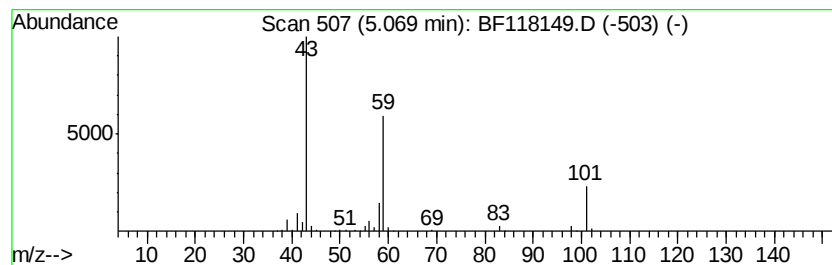
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	14.91 ng	445052	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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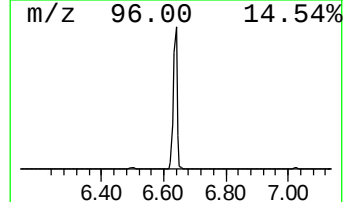
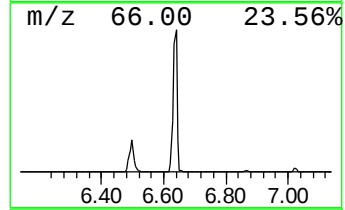
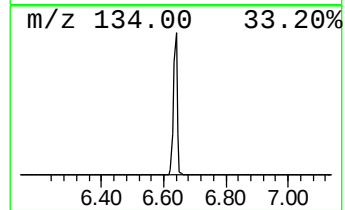
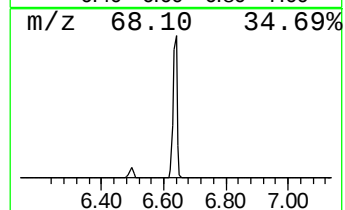
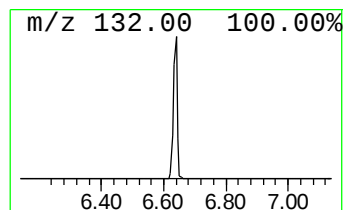
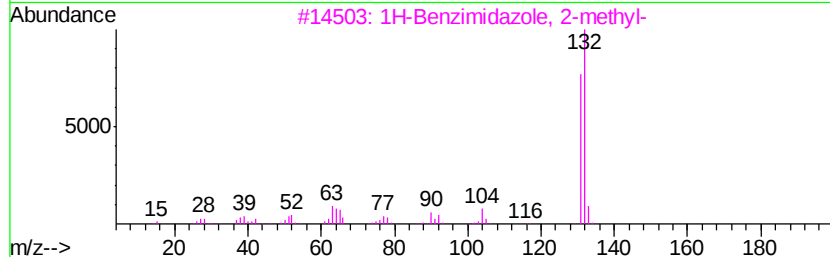
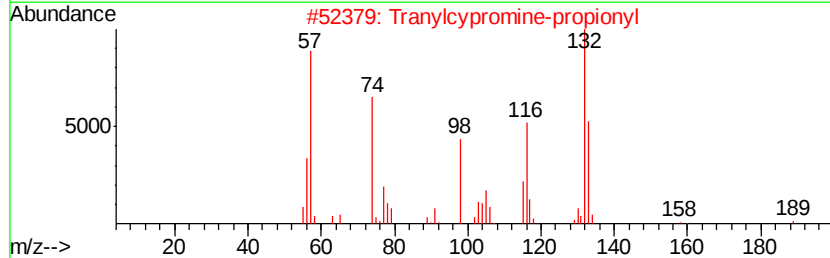
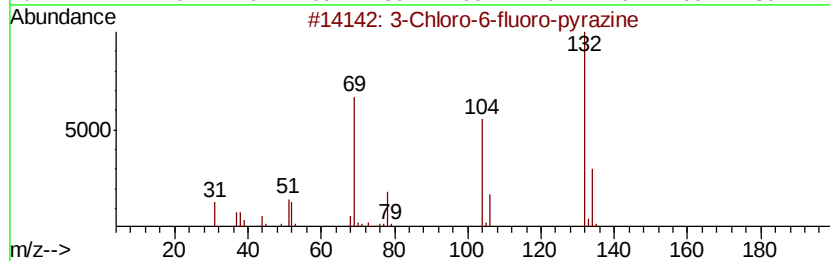
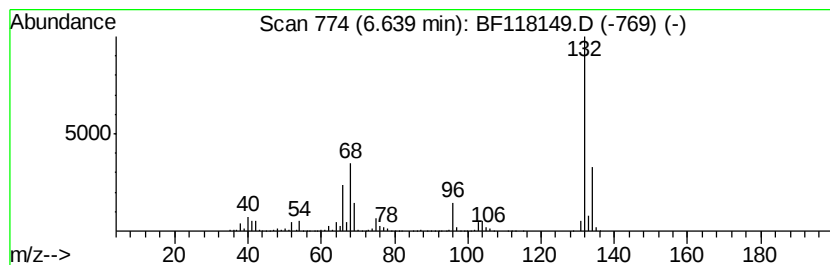
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown6.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	76.58 ng	2286360	1,4-Dichlorobenzene-d4	6.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	16
3			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
4			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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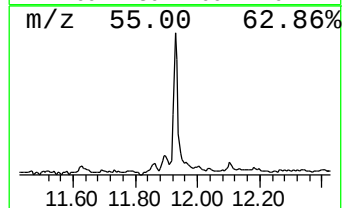
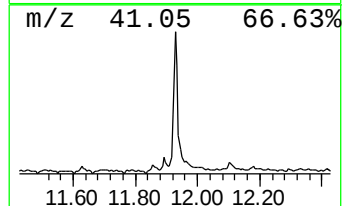
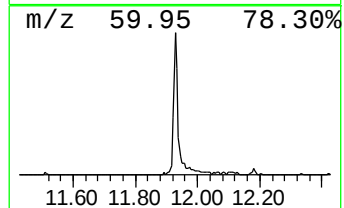
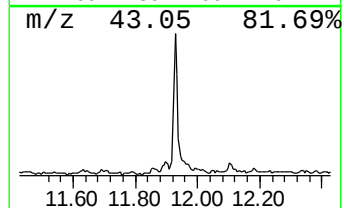
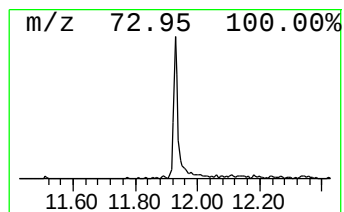
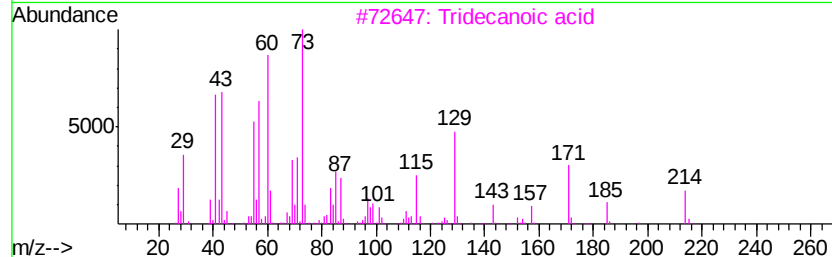
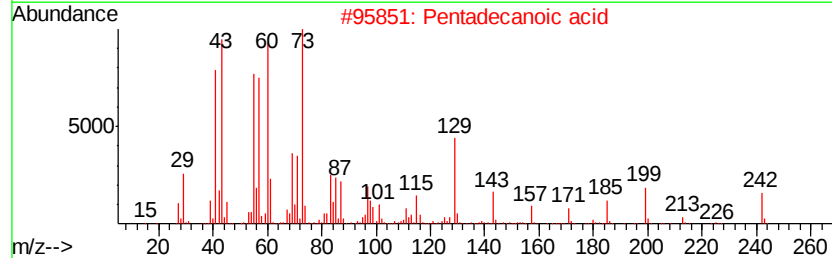
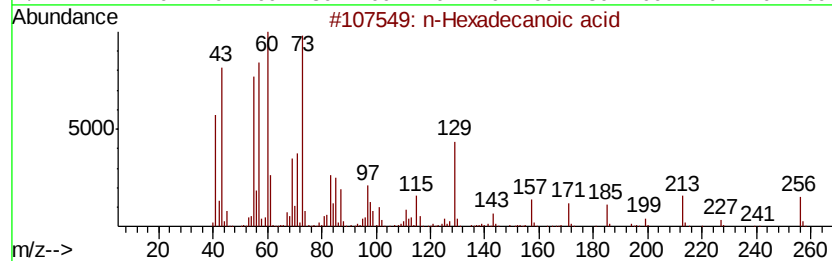
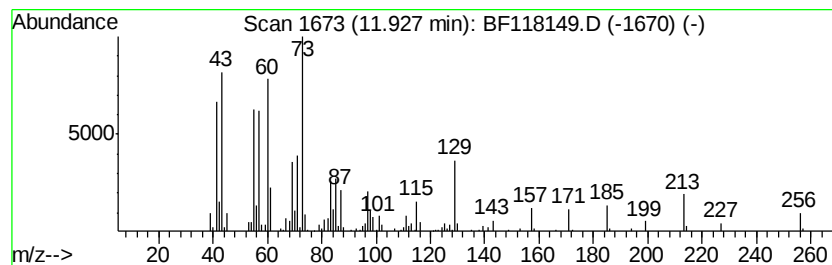
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	9.28 ng	363831	Phenanthrene-d10	11.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3		Tridecanoic acid	214	C13H26O2	000638-53-9	83
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5		Estra-1,3,5(10)-trien-17.beta.-ol	256	C18H24O	002529-64-8	25



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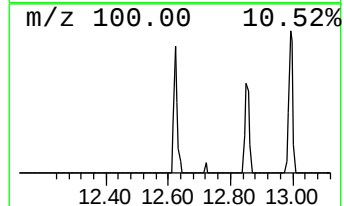
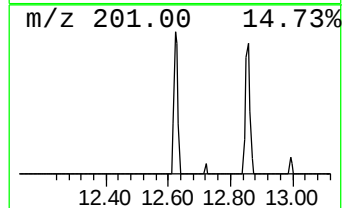
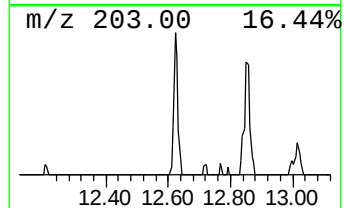
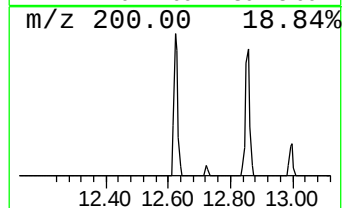
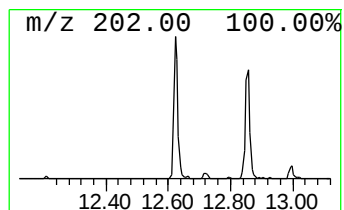
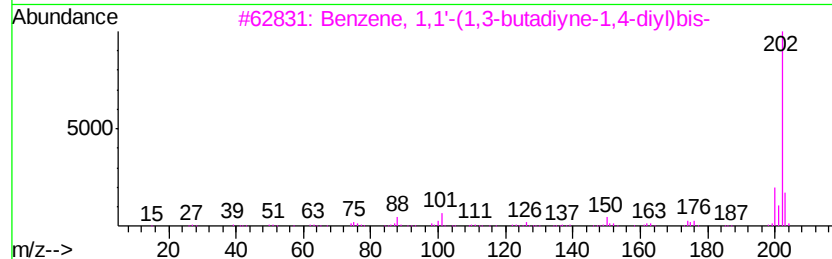
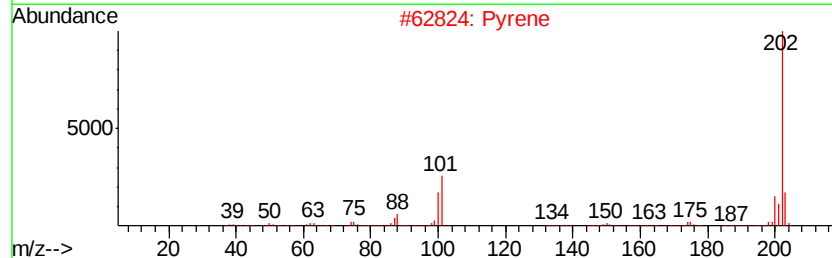
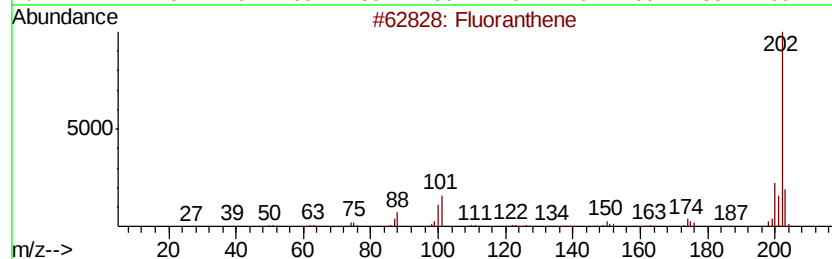
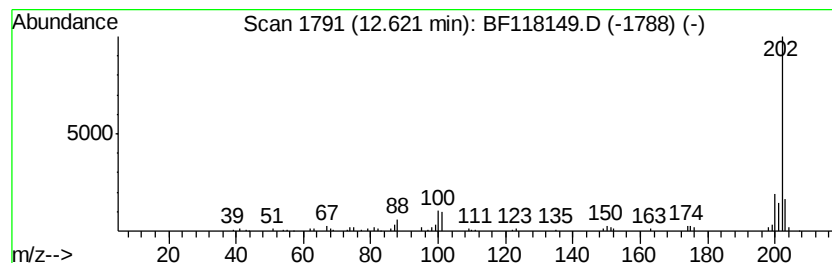
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,1'-(1,3-butadiyn... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.62	2.19 ng	85761	Phenanthrene-d10	11.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Fluoranthene	202	C16H10	000206-44-0	95
2		Pyrene	202	C16H10	000129-00-0	89
3		Benzene, 1,1'-(1,3-butadiyne-1,4...	202	C16H10	000886-66-8	81
4		4-Methoxy-naphthalene-1-carboxyl...	202	C12H10O3	1000317-85-4	40
5		Succinic acid, di(4-cyanophenyl)...	320	C18H12N2O4	1000360-71-2	28



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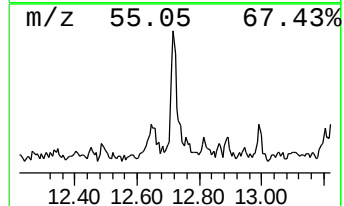
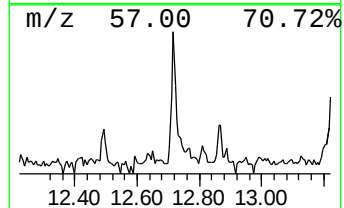
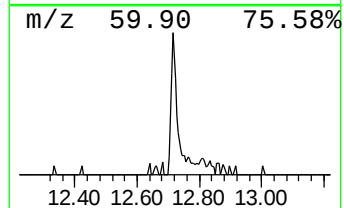
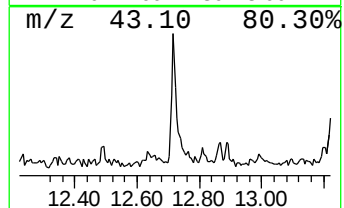
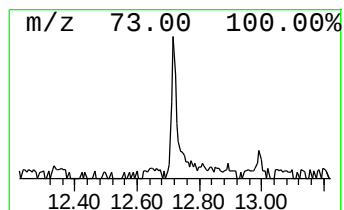
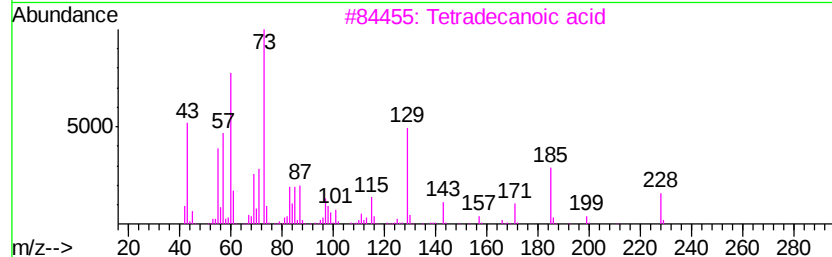
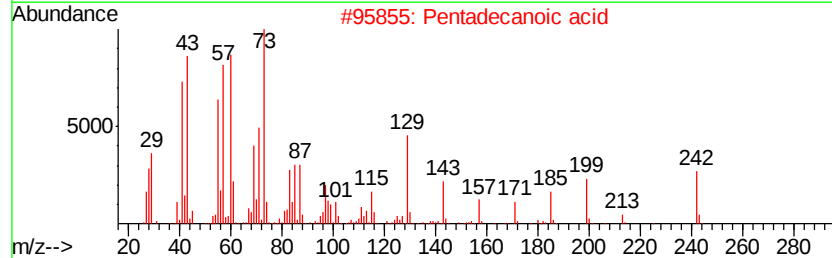
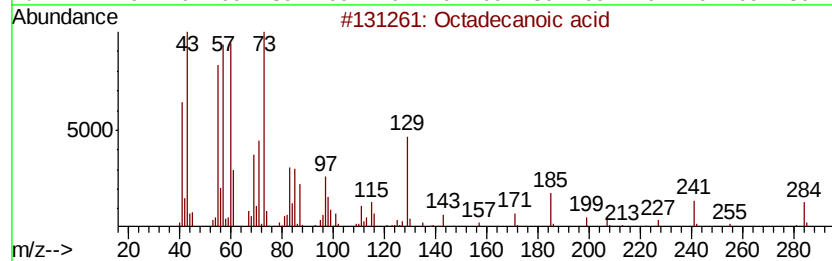
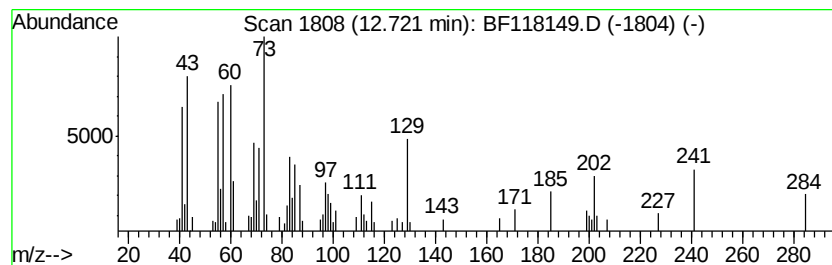
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.72	2.60 ng	101968	Phenanthrene-d10	11.40

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	81
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
4	n-Decanoic acid	172	C10H20O2	000334-48-5	53
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120919\
Data File : BF118149.D
Acq On : 9 Dec 2019 17:29
Operator : JU
Sample : K6187-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RPC-BOT-S

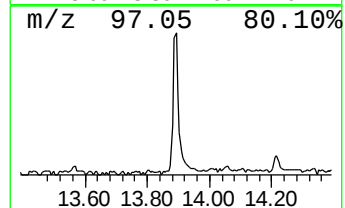
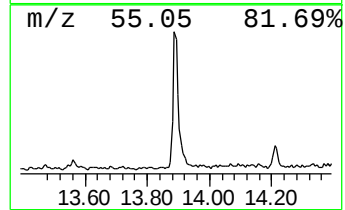
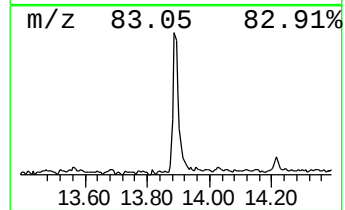
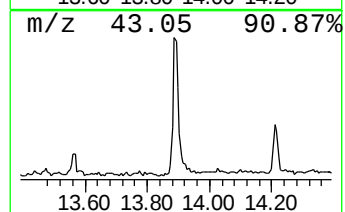
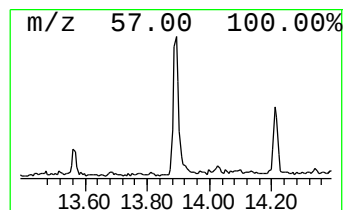
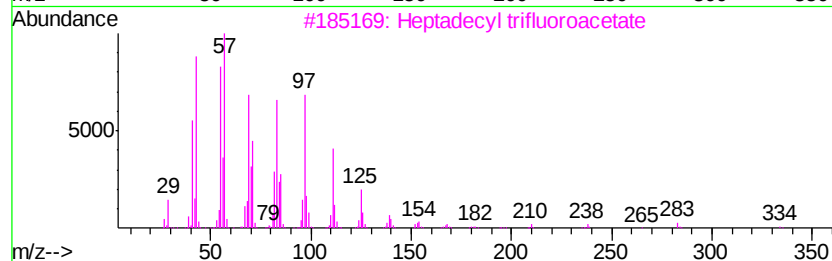
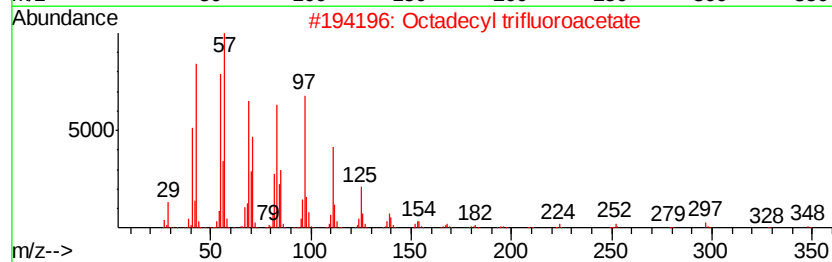
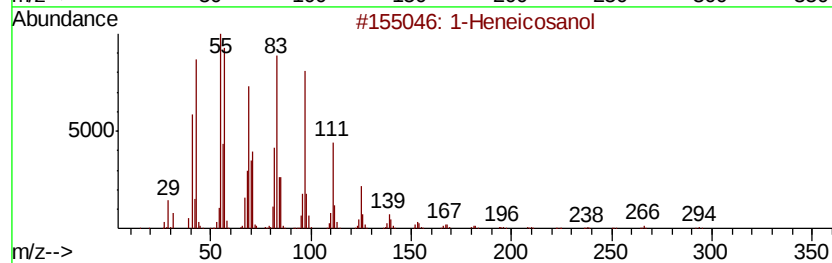
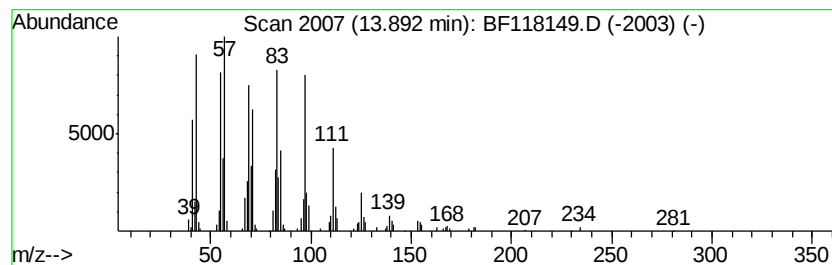
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	9.32 ng	308029	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosanol	312	C21H44O	015594-90-8	93
2		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	93
3		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	93
4		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	93
5		Behenic alcohol	326	C22H46O	000661-19-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120919\
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Acq On : 9 Dec 2019 17:29
Operator : JU
Sample : K6187-02
Misc :
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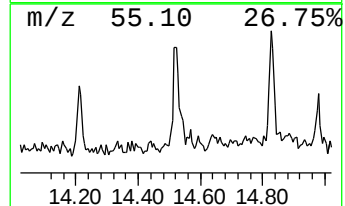
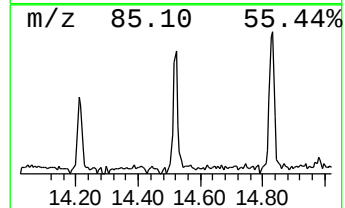
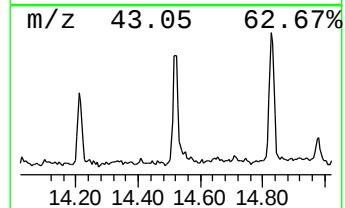
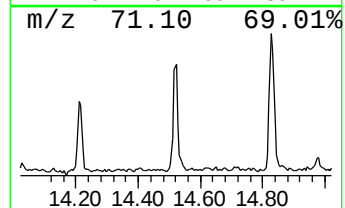
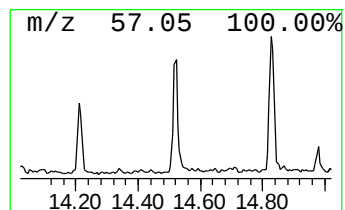
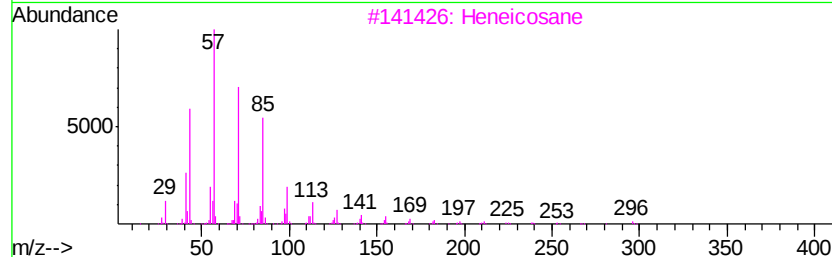
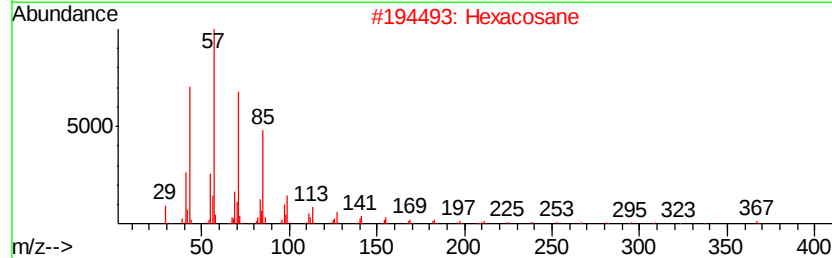
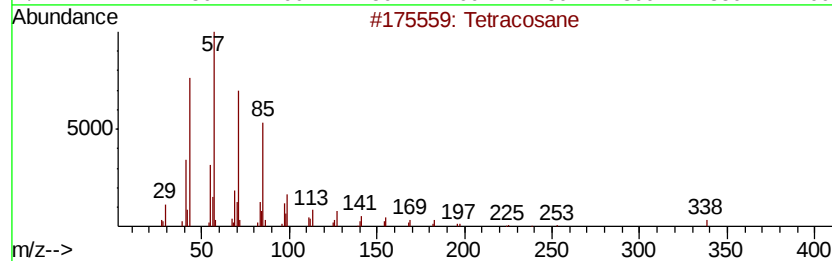
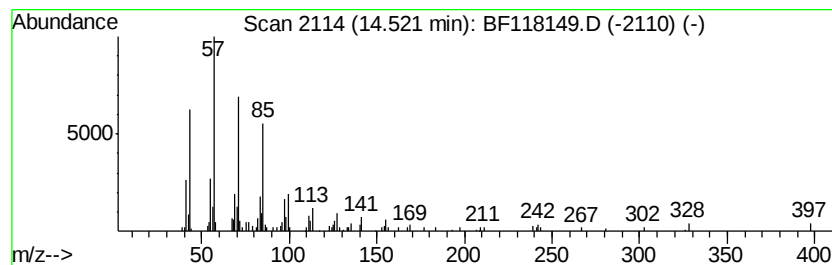
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetracosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	4.17 ng	137940	Chrysene-d12	14.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetracosane	338 C24H50	000646-31-1 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		Hentriacontane	437 C31H64	000630-04-6 91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120919\
Data File : BF118149.D
Acq On : 9 Dec 2019 17:29
Operator : JU
Sample : K6187-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

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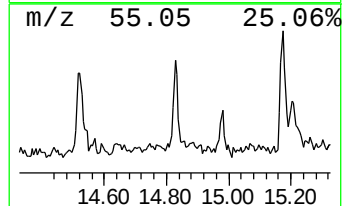
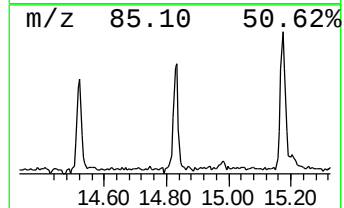
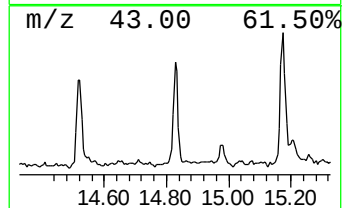
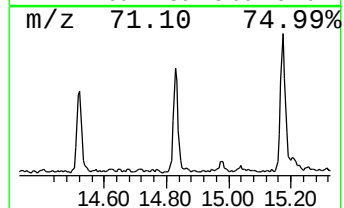
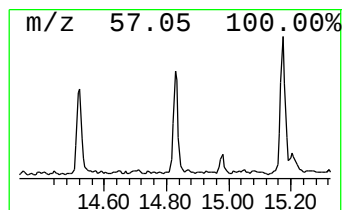
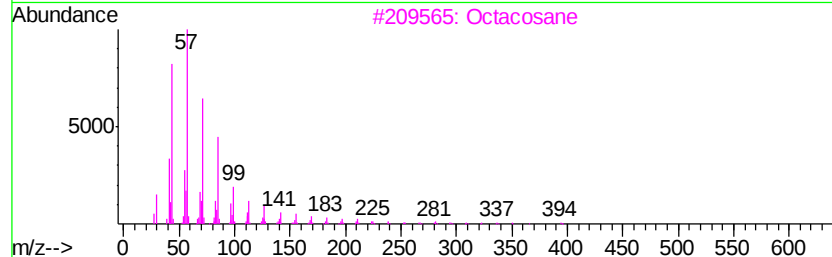
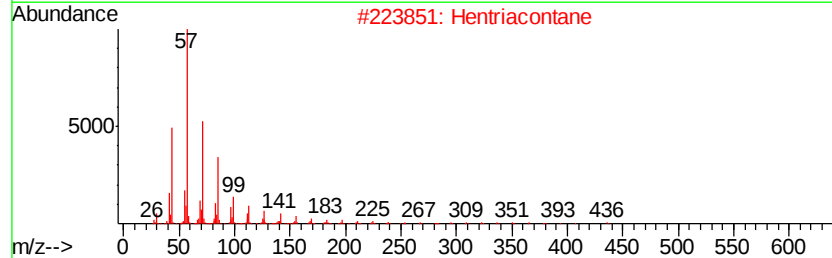
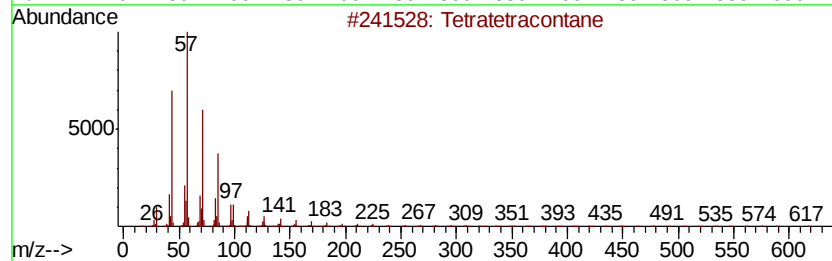
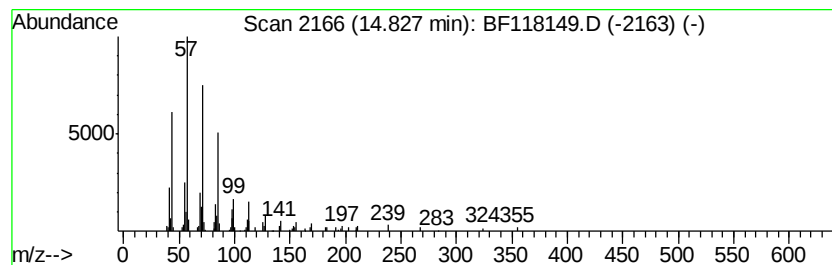
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hentriacontane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	2.53 ng	114831	Perylene-d12	15.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	91
2			Hentriacontane	437	C31H64	000630-04-6	91
3			Octacosane	394	C28H58	000630-02-4	91
4			Hexacosane	366	C26H54	000630-01-3	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120919\
Data File : BF118149.D
Acq On : 9 Dec 2019 17:29
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Misc :
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Instrument :
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ClientSampleId :
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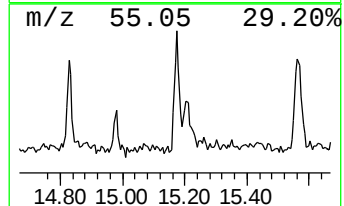
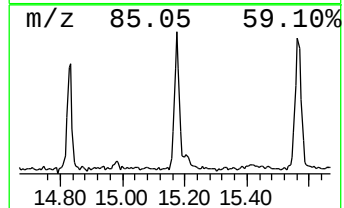
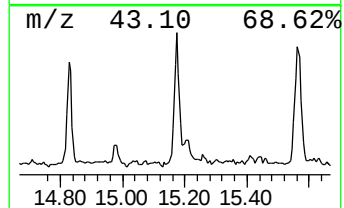
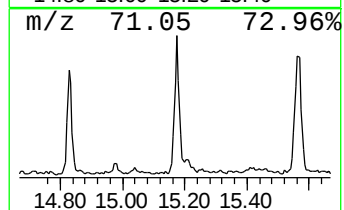
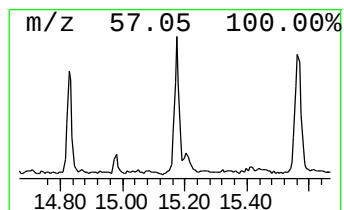
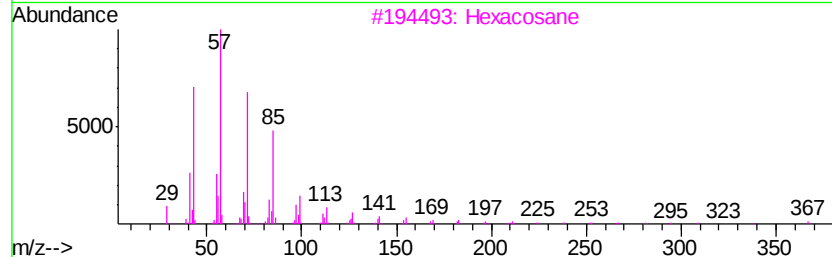
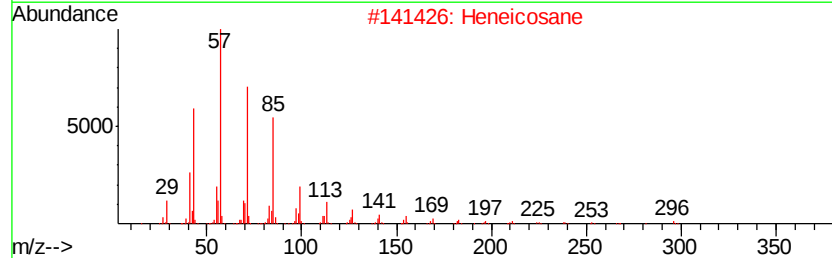
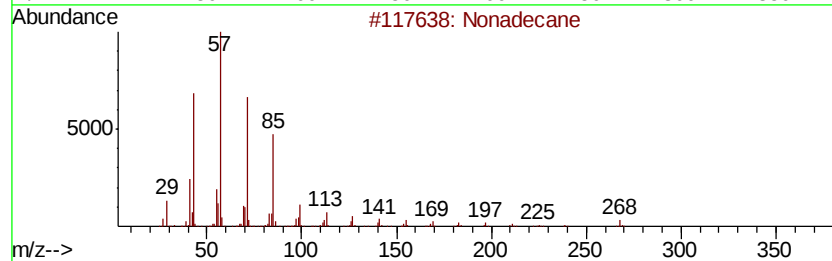
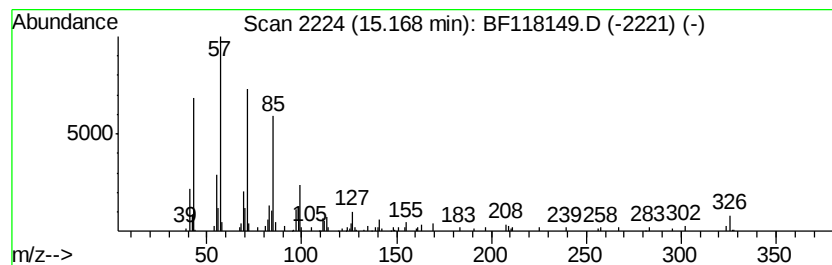
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Nonadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	3.85 ng	174741	Perylene-d12	15.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	95
2		Heneicosane	296	C21H44	000629-94-7	90
3		Hexacosane	366	C26H54	000630-01-3	90
4		Docosane	310	C22H46	000629-97-0	90
5		Triacontane	422	C30H62	000638-68-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120919\
Data File : BF118149.D
Acq On : 9 Dec 2019 17:29
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Misc :
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ClientSampleId :
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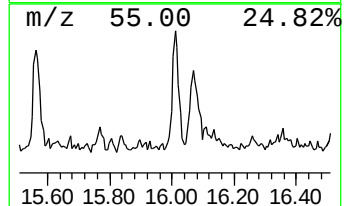
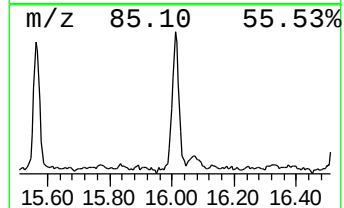
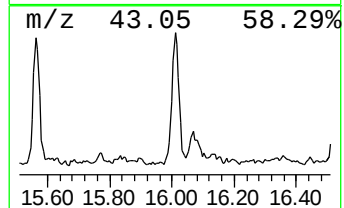
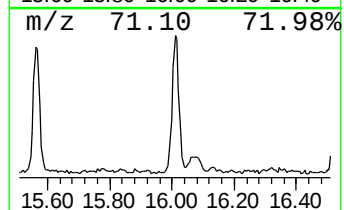
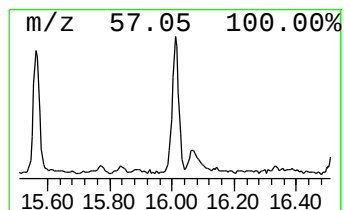
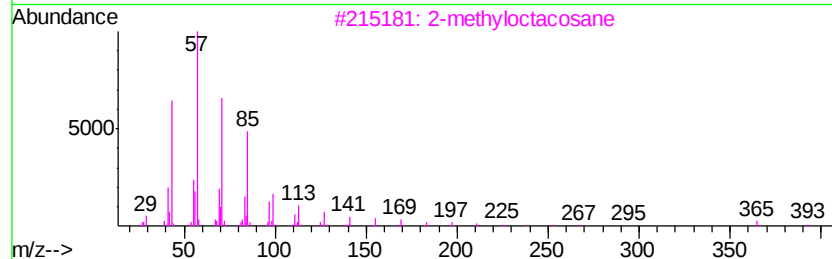
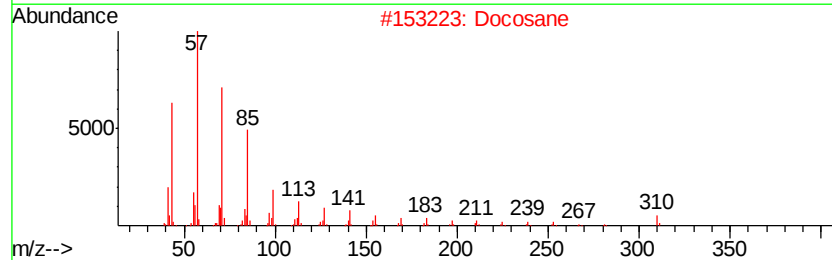
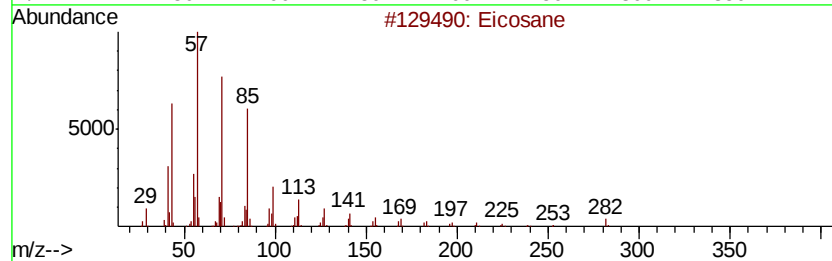
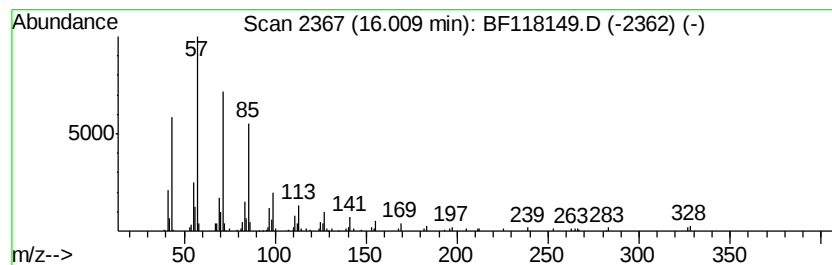
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Eicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.01	4.64 ng	210780	Perylene-d12	15.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Docosane	310	C22H46	000629-97-0	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Pentacosane	352	C25H52	000629-99-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120919\
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Acq On : 9 Dec 2019 17:29
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ClientSampleId :
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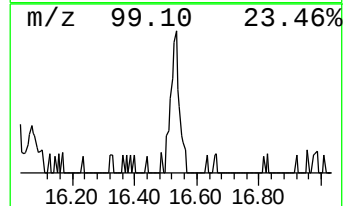
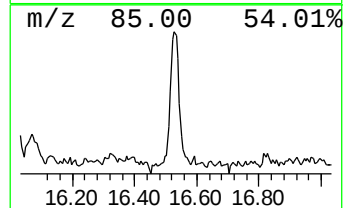
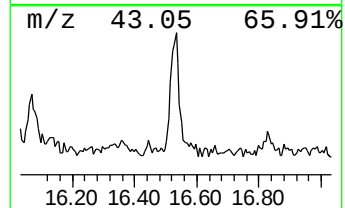
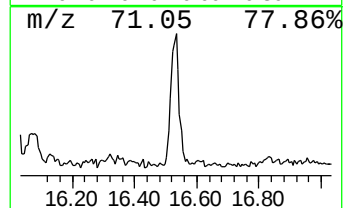
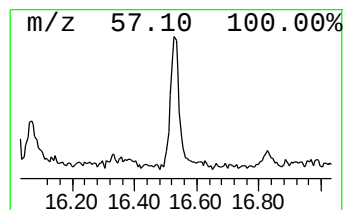
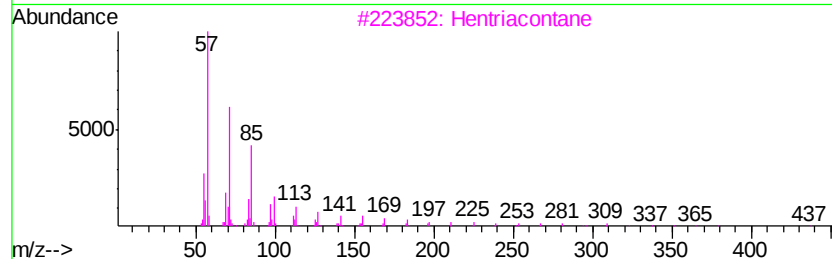
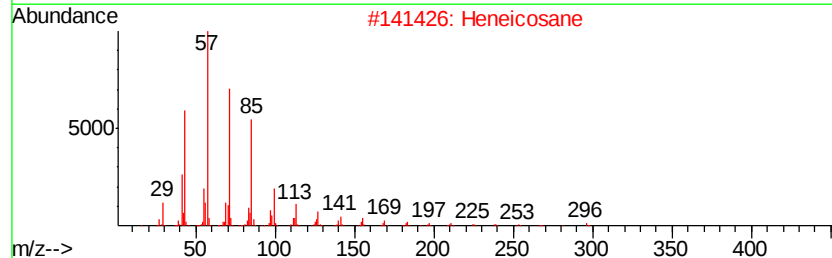
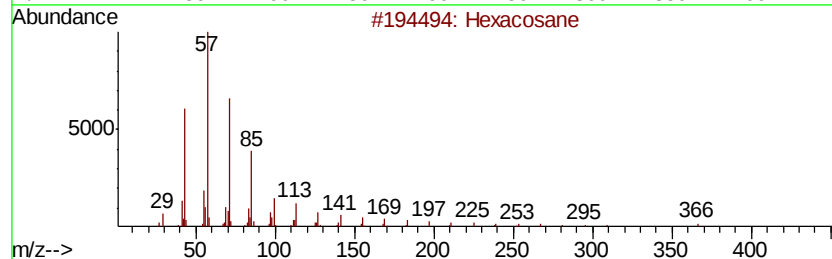
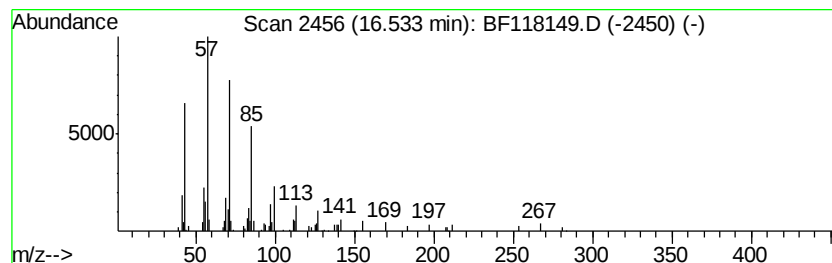
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.53	2.52 ng	114670	Perylene-d12	15.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Hentriacontane	437	C31H64	000630-04-6	91
4		Triacontane	422	C30H62	000638-68-6	91
5		Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF120919\
Data File : BF118149.D
Acq On : 9 Dec 2019 17:29
Operator : JU
Sample : K6187-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
RPC-BOT-S

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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
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unknown6.64	6.64	76.6	ng	2286360	1	6.87	597144	20.0
n-Hexadecanoic acid	11.93	9.3	ng	363831	4	11.40	784231	20.0
Benzene, 1,1'-(1,...	12.62	2.2	ng	85761	4	11.40	784231	20.0
Octadecanoic acid	12.72	2.6	ng	101968	4	11.40	784231	20.0
1-Heneicosanol	13.89	9.3	ng	308029	5	14.06	660946	20.0
Tetracosane	14.52	4.2	ng	137940	5	14.06	660946	20.0
Hentriacontane	14.83	2.5	ng	114831	6	15.55	908724	20.0
Nonadecane	15.17	3.9	ng	174741	6	15.55	908724	20.0
Eicosane	16.01	4.6	ng	210780	6	15.55	908724	20.0
Hexacosane	16.53	2.5	ng	114670	6	15.55	908724	20.0