

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031122\
 Data File : BP009377.D
 Acq On : 11 Mar 2022 22:01
 Operator : CG/JU
 Sample : N1886-01 20X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 2822-D

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP009377.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.893	148	154	177	rBV3	102363	559026	3.85%	0.869%
2	5.428	409	415	422	rBV	133127	228257	1.57%	0.355%
3	7.016	674	685	698	rBV	118798	264993	1.83%	0.412%
4	7.816	815	821	830	rVB	1290467	2163660	14.91%	3.364%
5	9.058	1023	1032	1041	rVB	561291	1010522	6.96%	1.571%
6	10.610	1288	1296	1315	rVB	1878528	3453549	23.80%	5.369%
7	11.534	1445	1453	1467	rBV4	101956	426540	2.94%	0.663%
8	13.081	1709	1716	1725	rBV	265244	469063	3.23%	0.729%
9	14.122	1887	1893	1906	rBV3	68615	213961	1.47%	0.333%
10	14.457	1944	1950	1972	rBV	2319615	4142162	28.54%	6.439%
11	15.469	2117	2122	2133	rBV6	163093	386618	2.66%	0.601%
12	15.651	2147	2153	2157	rBV2	224286	390643	2.69%	0.607%
13	15.698	2157	2161	2164	rBV	163904	233822	1.61%	0.363%
14	15.957	2200	2205	2211	rBV	213173	382437	2.64%	0.595%
15	16.198	2241	2246	2253	rBV2	150646	308582	2.13%	0.480%
16	16.269	2253	2258	2268	rVB6	102489	258547	1.78%	0.402%
17	16.998	2377	2382	2386	rBV	320088	445436	3.07%	0.692%
18	17.222	2414	2420	2430	rVB	3076982	4701925	32.40%	7.309%
19	17.710	2499	2503	2505	rBV2	204005	292736	2.02%	0.455%
20	17.739	2505	2508	2514	rVB2	236152	348826	2.40%	0.542%
21	18.128	2570	2574	2580	rBV2	493578	833929	5.75%	1.296%
22	18.234	2590	2592	2600	rVB5	145558	221192	1.52%	0.344%
23	18.416	2619	2623	2629	rVB	378979	550542	3.79%	0.856%
24	18.863	2696	2699	2704	rVB2	359754	503384	3.47%	0.783%
25	19.028	2724	2727	2731	rBV	451762	517347	3.56%	0.804%
26	19.551	2811	2816	2821	rBV	10532075	13182341	90.83%	20.493%
27	19.592	2821	2823	2827	rVB	502360	547320	3.77%	0.851%
28	20.675	3003	3007	3012	rBV3	2645240	4902509	33.78%	7.621%
29	20.928	3047	3050	3056	rBV	2541544	3731823	25.71%	5.801%
30	21.333	3116	3119	3123	rBV	3187439	4142758	28.55%	6.440%
31	21.451	3135	3139	3144	rBV	9559251	14512707	100.00%	22.561%

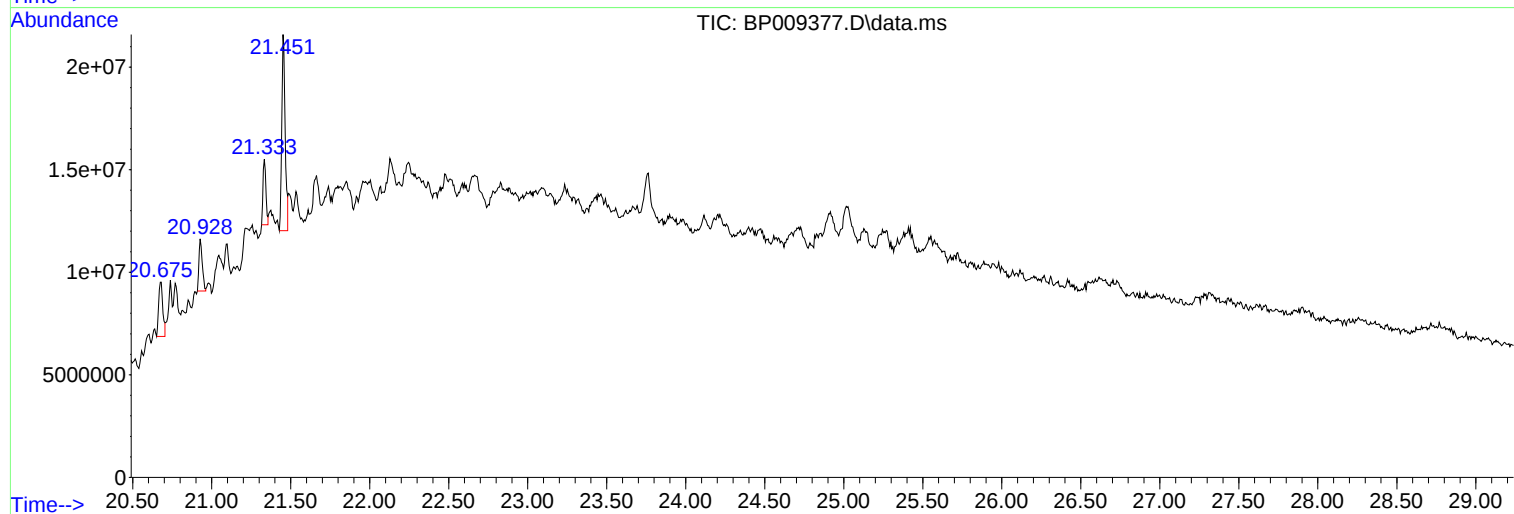
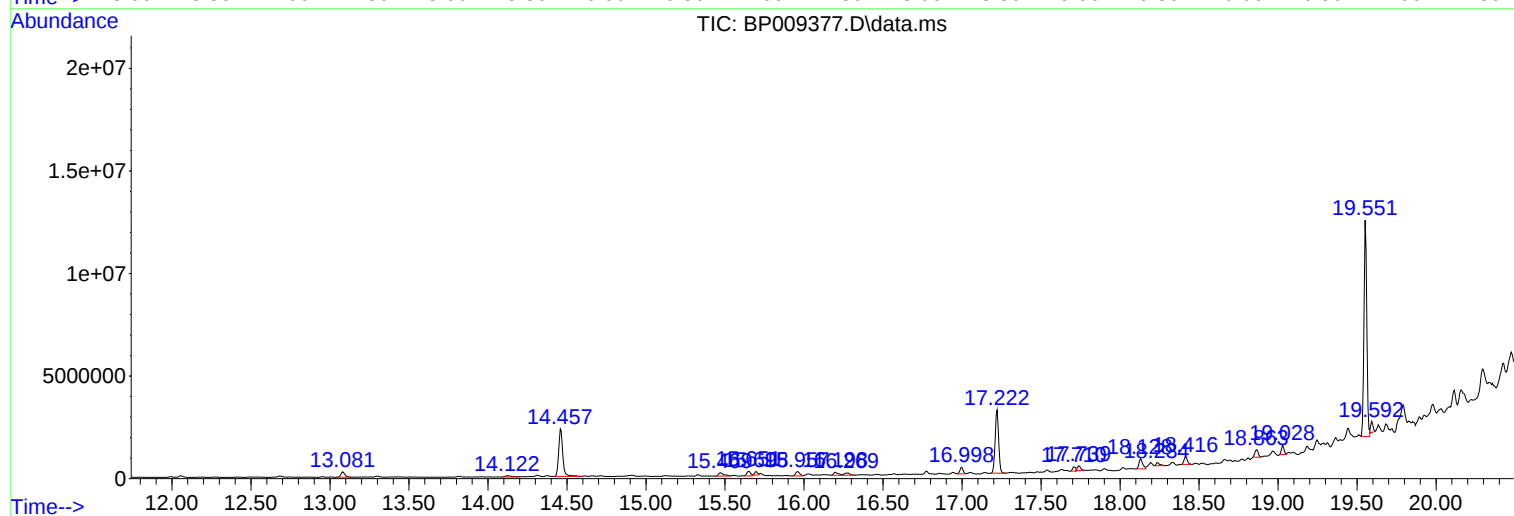
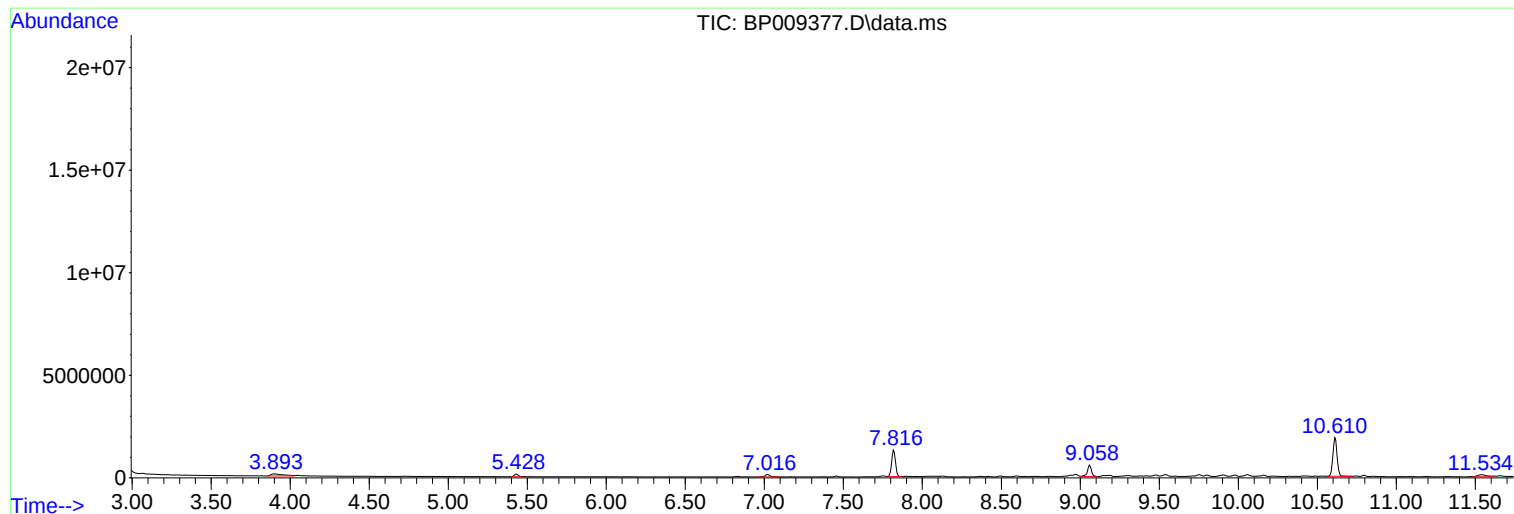
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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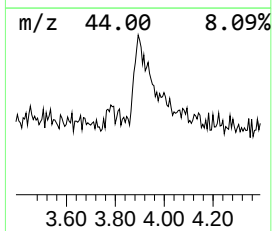
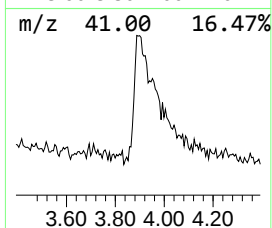
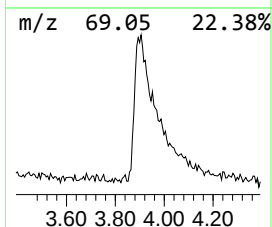
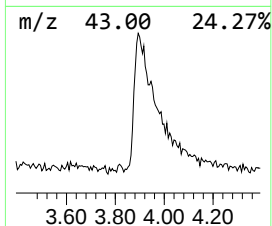
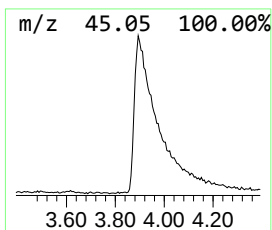
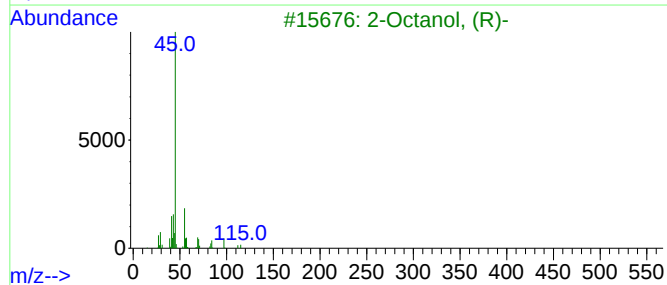
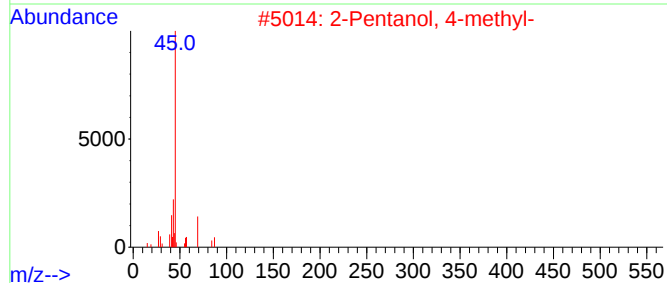
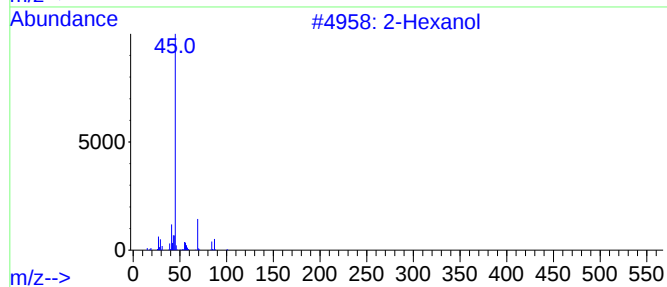
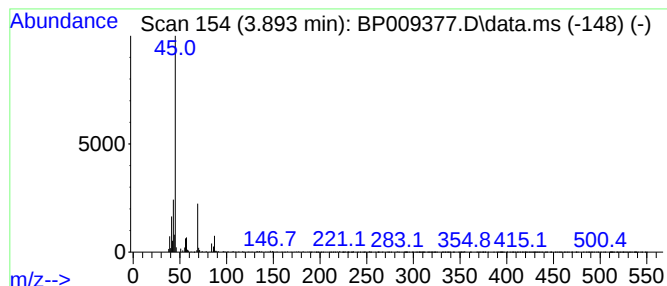
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Hexanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.893	5.17 ng	559026	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol			102	C6H14O	000626-93-7	83
2	2-Pentanol, 4-methyl-			102	C6H14O	000108-11-2	83
3	2-Octanol, (R)-			130	C8H18O	005978-70-1	78
4	2-Octanol, (S)-			130	C8H18O	006169-06-8	64
5	2-Octanol			130	C8H18O	000123-96-6	64



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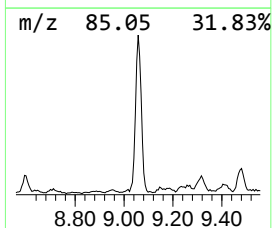
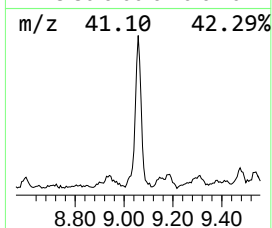
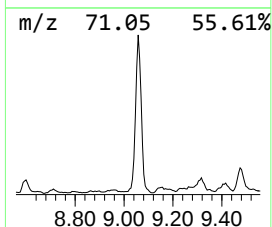
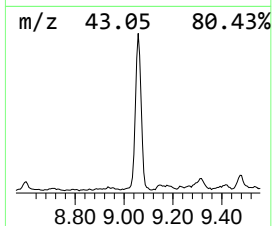
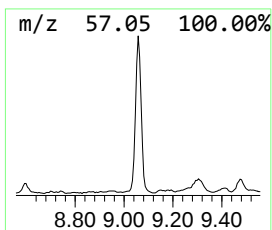
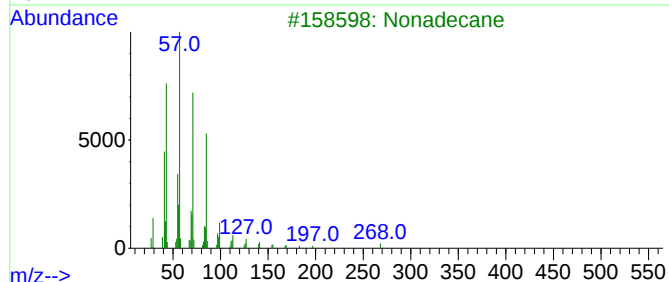
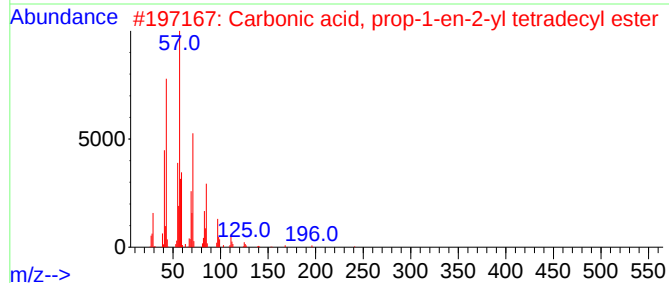
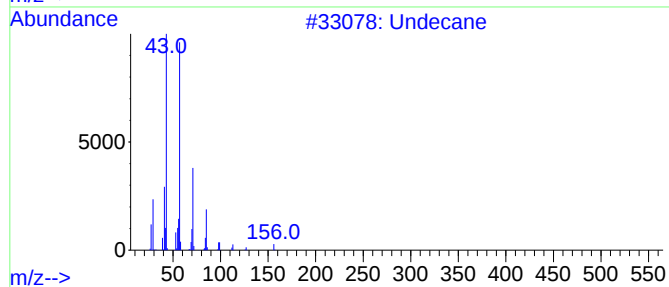
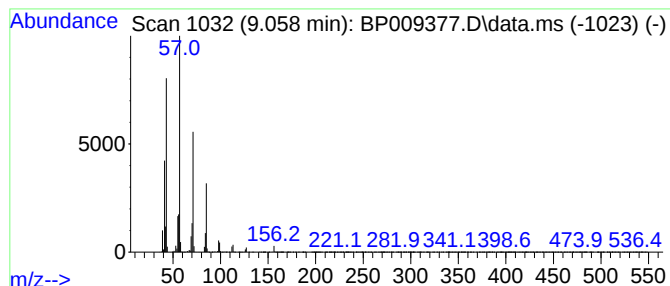
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Undecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.058	9.34 ng	1010520	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	91
2	Carbonic acid, prop-1-en-2-yl te...			298	C18H34O3	1000382-90-9	90
3	Nonadecane			268	C19H40	000629-92-5	87
4	Tridecane			184	C13H28	000629-50-5	86
5	Carbonic acid, prop-1-en-2-yl tr...			284	C17H32O3	1000382-90-8	86



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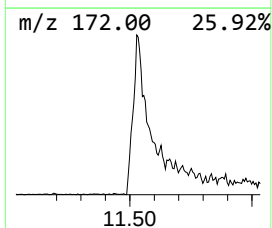
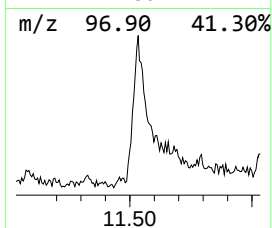
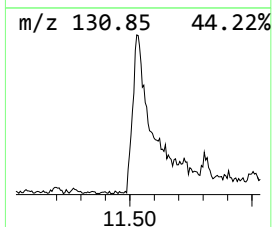
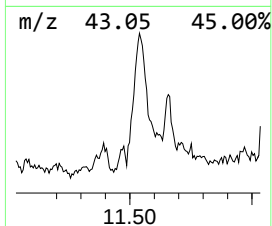
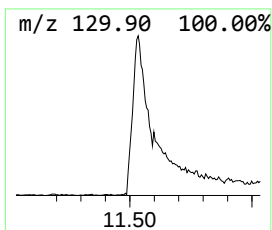
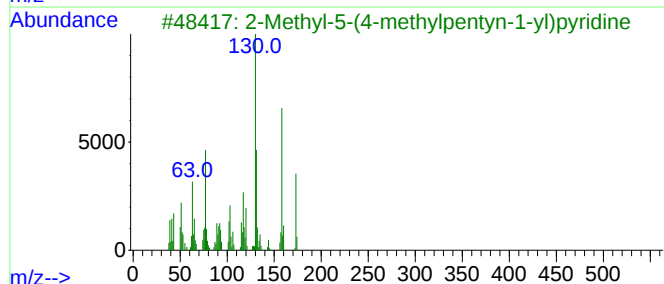
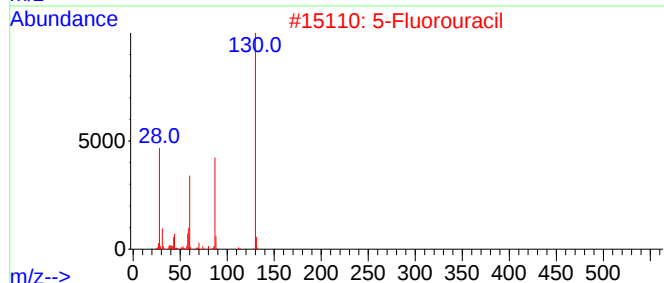
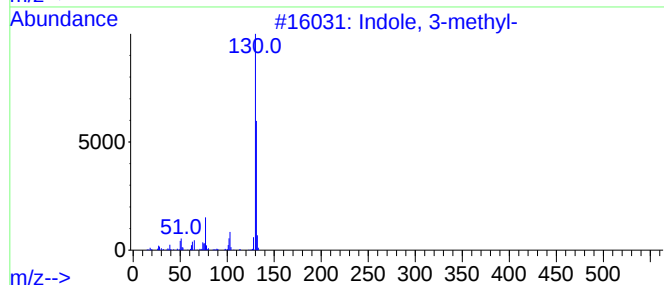
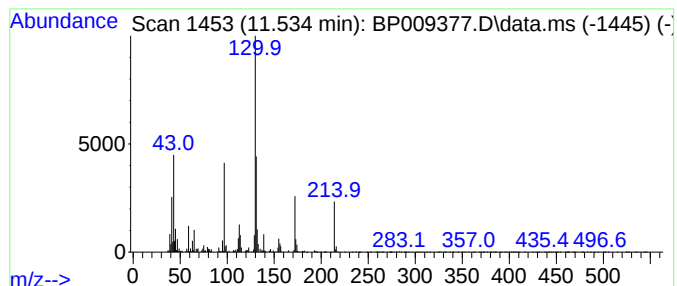
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 unknown11.534 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.534	2.47 ng	426540	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indole, 3-methyl-	131	C9H9N	000083-34-1	37
2			5-Fluorouracil	130	C4H3FN2O2	000051-21-8	27
3			2-Methyl-5-(4-methylpentyn-1-yl)...	173	C12H15N	016232-31-8	25
4			4-Fluoroimidazole-5-carboxylic a...	172	C7H9FN2O2	1000116-94-4	25
5			2'-Acetyl-1'-(1,2-dihydroquinox...	303	C19H17N3O	1000103-60-8	25



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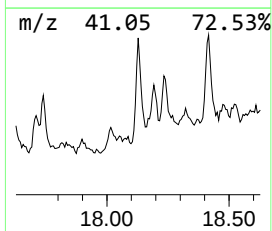
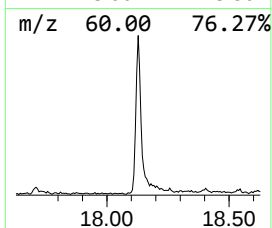
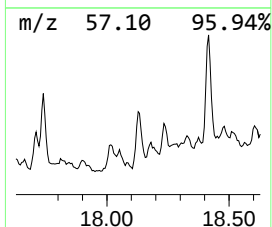
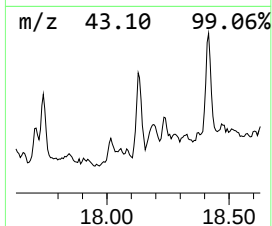
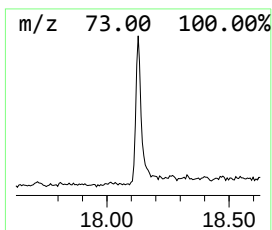
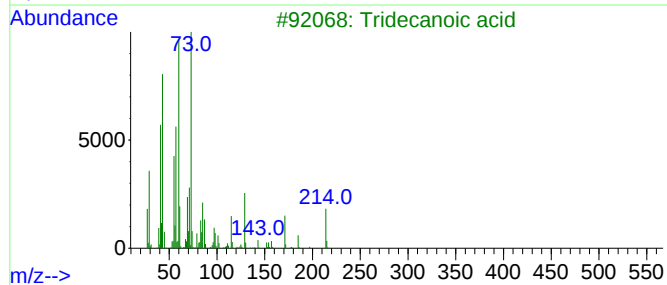
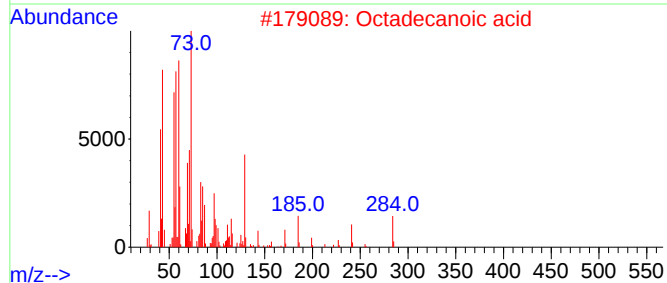
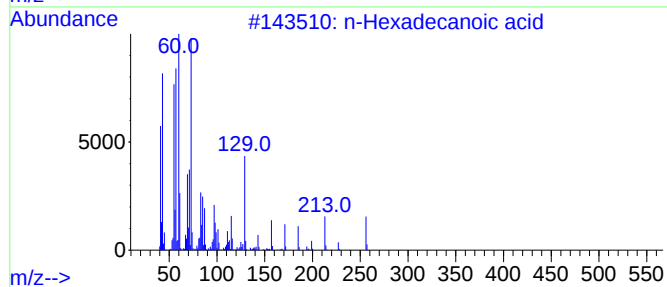
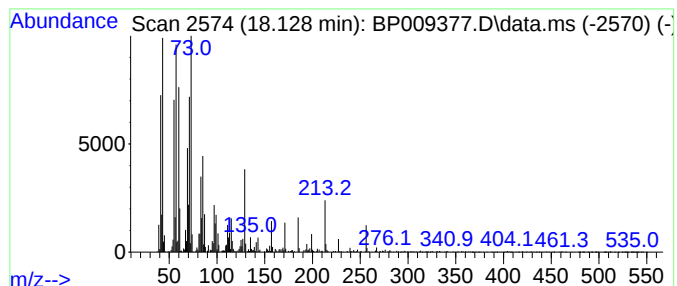
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.128	3.55 ng	833929	Phenanthrene-d10	17.222

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	68
3			Tridecanoic acid	214	C13H26O2	000638-53-9	64
4			l-(+)-Ascorbic acid 2,6-dihexade...	652	C38H68O8	028474-90-0	64
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	58



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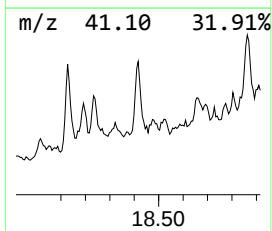
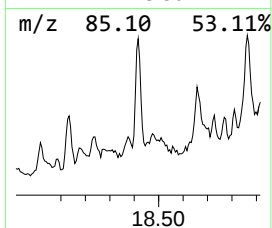
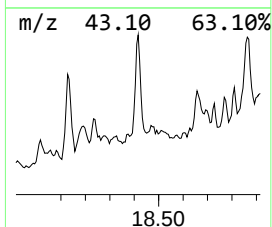
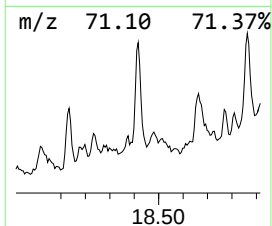
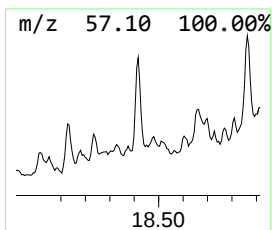
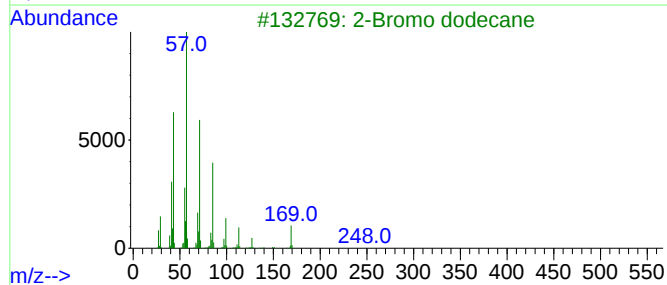
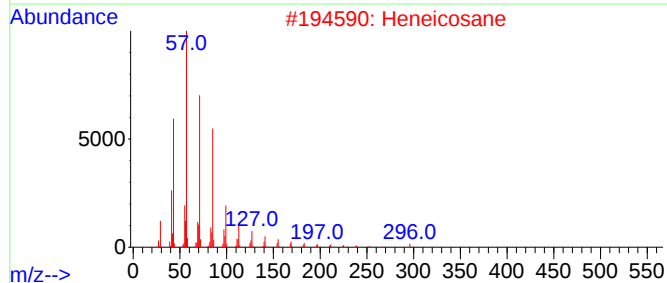
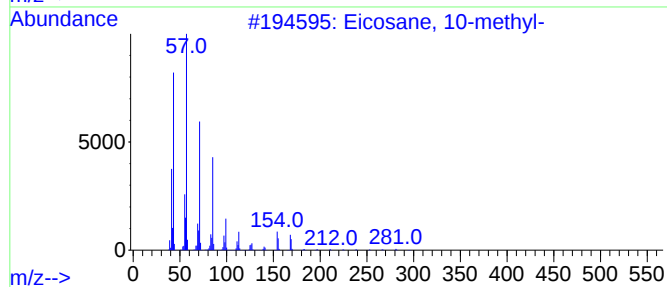
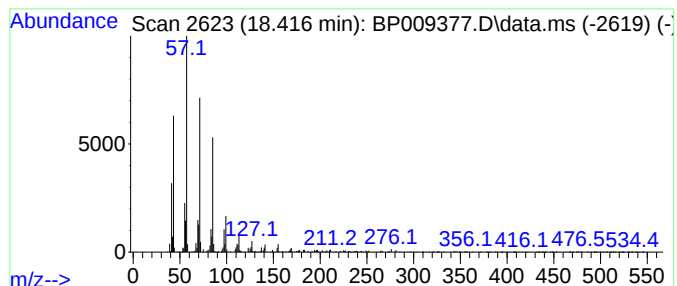
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Peak Number 5 Eicosane, 10-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.416	2.34 ng	550542	Phenanthrene-d10	17.222

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
2			Heneicosane	296	C21H44	000629-94-7	91
3			2-Bromo dodecane	248	C12H25Br	013187-99-0	90
4			Heptadecane	240	C17H36	000629-78-7	87
5			Octadecane	254	C18H38	000593-45-3	83



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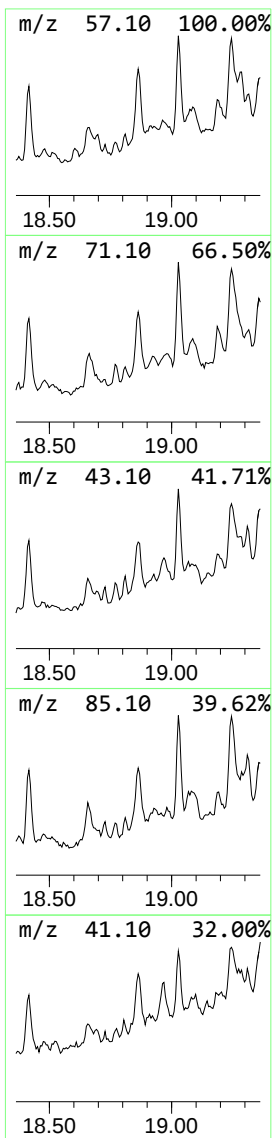
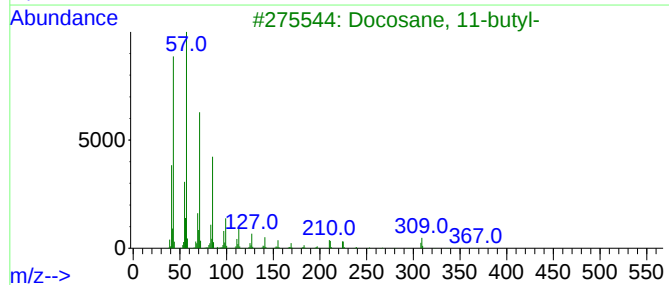
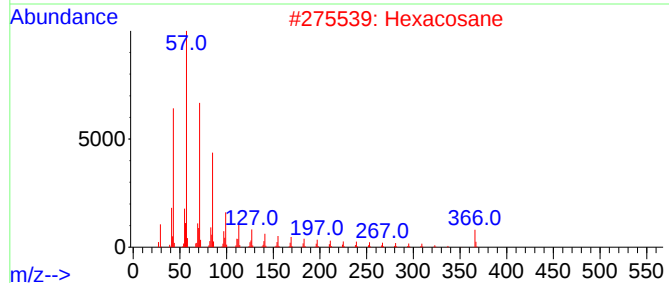
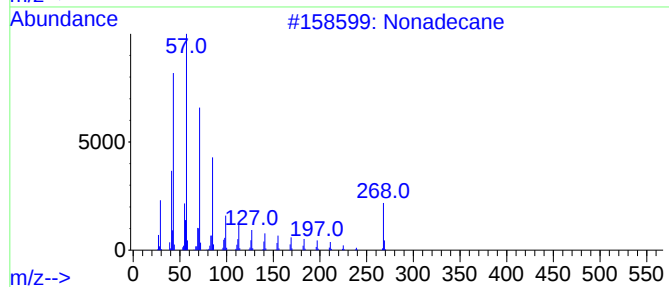
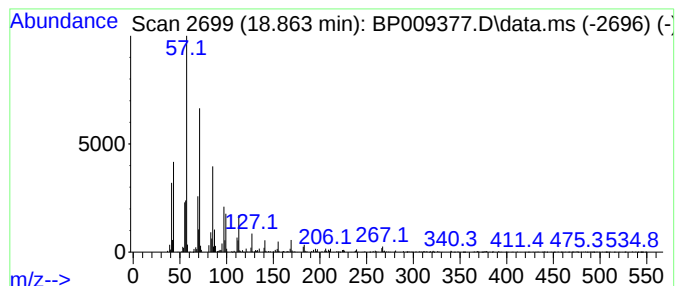
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.863	2.14 ng	503384	Phenanthrene-d10	17.222	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	83
2	Hexacosane	366	C26H54	000630-01-3	83
3	Docosane, 11-butyl-	366	C26H54	013475-76-8	76
4	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	72
5	2-Methylhexacosane	380	C27H56	001561-02-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031122\
Data File : BP009377.D
Acq On : 11 Mar 2022 22:01
Operator : CG/JU
Sample : N1886-01 20X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
2822-D

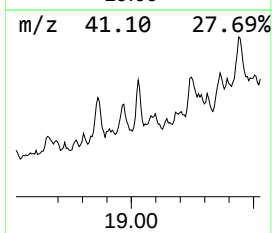
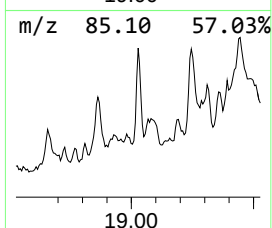
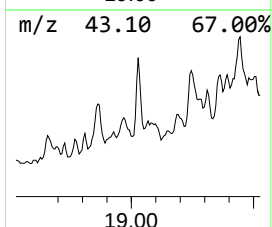
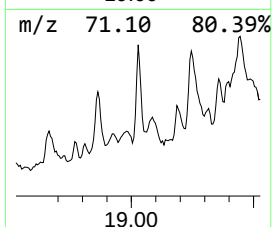
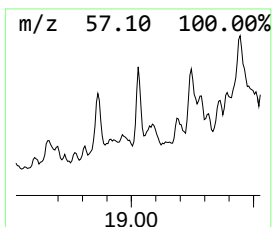
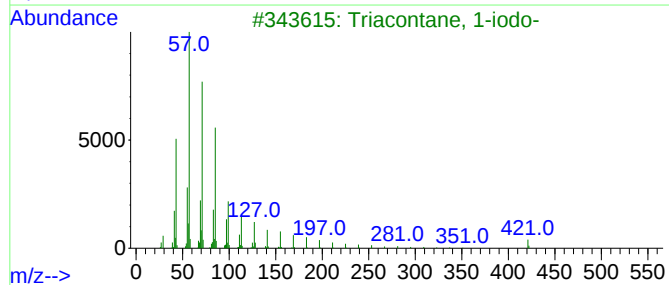
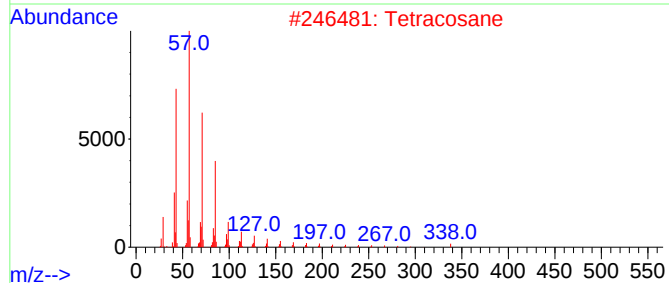
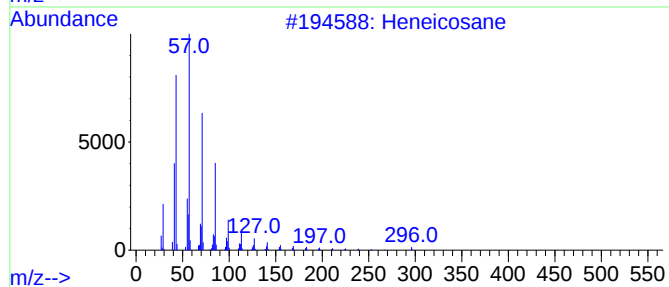
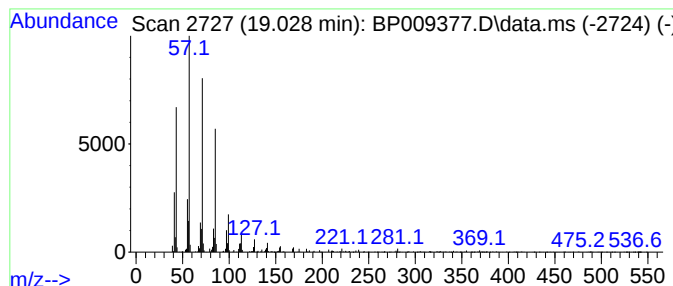
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Heneicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.028	2.20 ng	517347	Phenanthrene-d10	17.222

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Tetracosane	338	C24H50	000646-31-1	91
3			Triacontane, 1-iodo-	548	C30H61I	1000406-32-3	90
4			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	87
5			Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031122\
Data File : BP009377.D
Acq On : 11 Mar 2022 22:01
Operator : CG/JU
Sample : N1886-01 20X
Misc :
ALS Vial : 20 Sample Multiplier: 1

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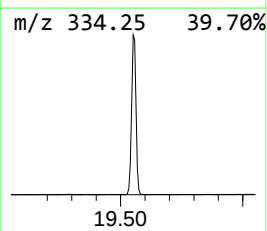
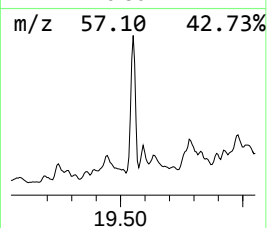
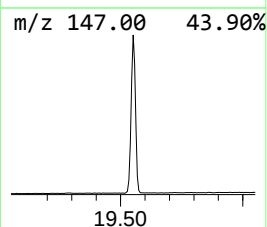
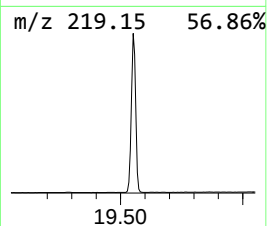
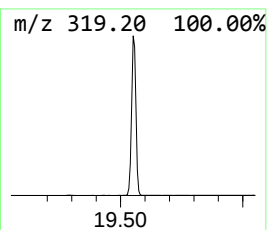
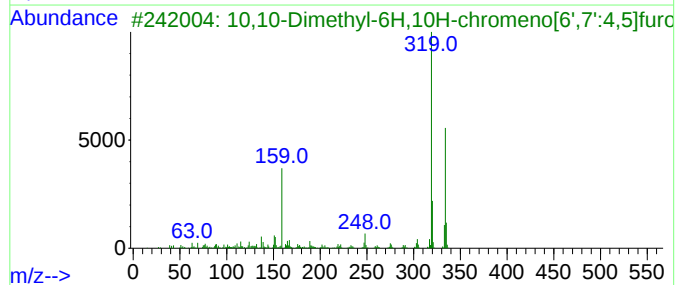
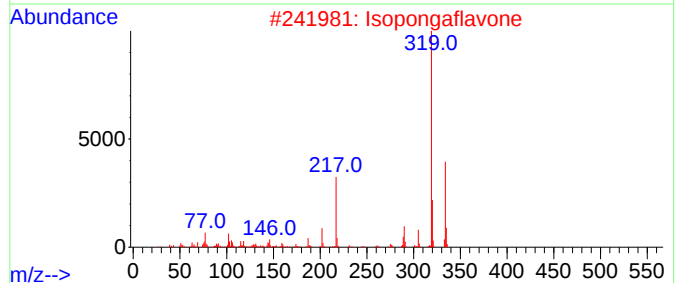
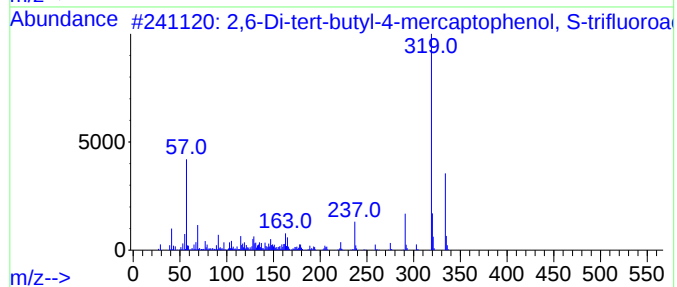
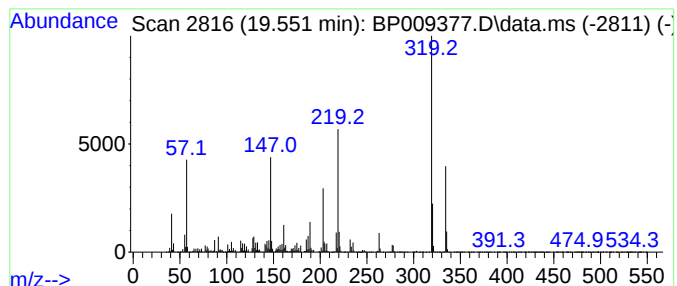
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown19.551 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.551	63.64 ng	13182300	Chrysene-d12	21.333

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6-Di-tert-butyl-4-mercaptophen...	334	C16H21F3O2S	1010467-17-3	46		
2	Isopongaflavone	334	C21H18O4	064125-33-3	46		
3	10,10-Dimethyl-6H,10H-chromeno[6...	334	C21H18O4	041347-50-6	43		
4	1H,6H-5,11b-Ethano[1,3]dioxolo[4...	319	C17H21NO5	003660-65-9	27		
5	Methyl 6-(naphthalen-2-yl)-5-oxo...	319	C20H17NO3	1000482-02-2	25		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031122\
Data File : BP009377.D
Acq On : 11 Mar 2022 22:01
Operator : CG/JU
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Misc :
ALS Vial : 20 Sample Multiplier: 1

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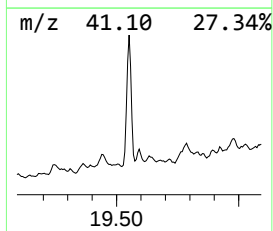
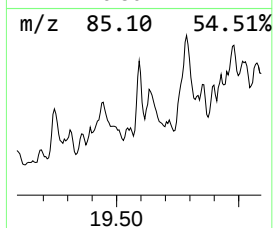
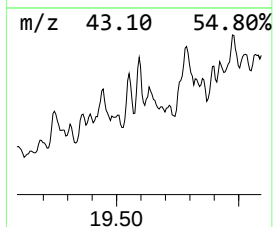
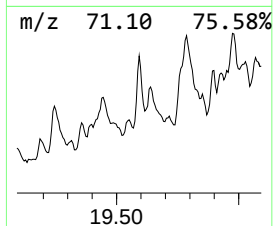
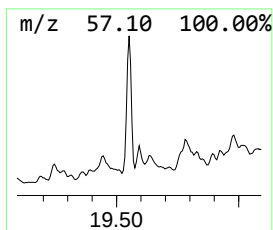
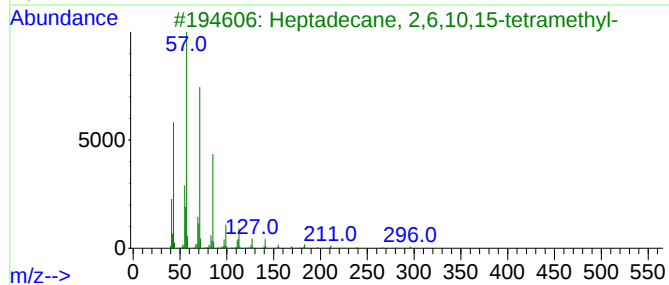
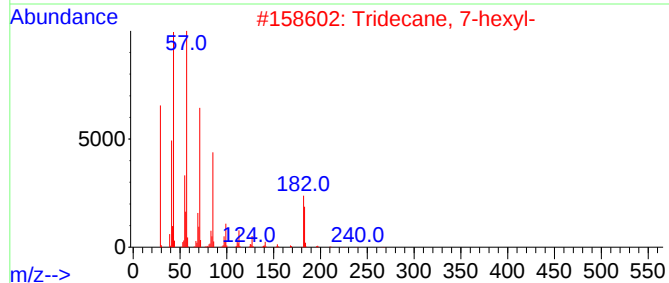
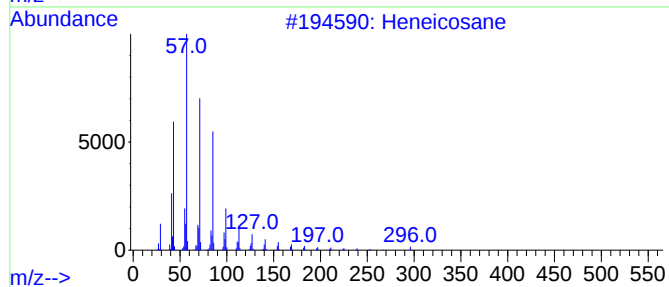
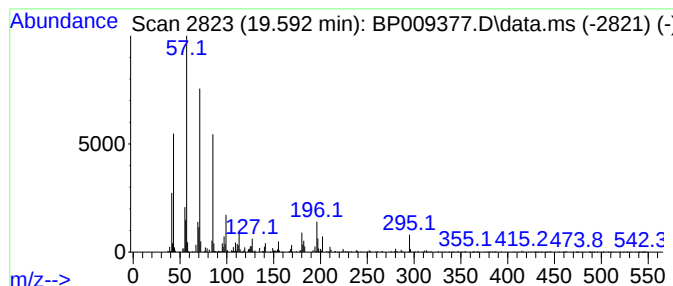
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Tridecane, 7-hexyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.592	2.64 ng	547320	Chrysene-d12	21.333

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	99
2		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	91
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5		Tetracosane	338	C24H50	000646-31-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031122\
Data File : BP009377.D
Acq On : 11 Mar 2022 22:01
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Misc :
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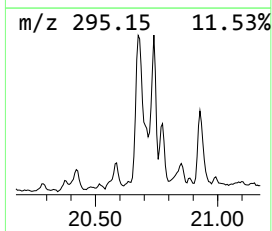
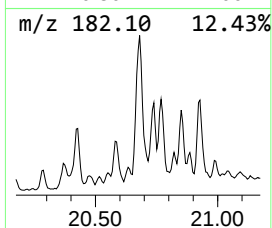
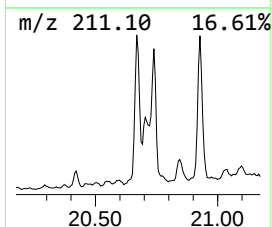
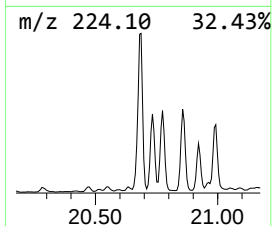
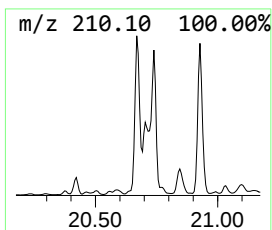
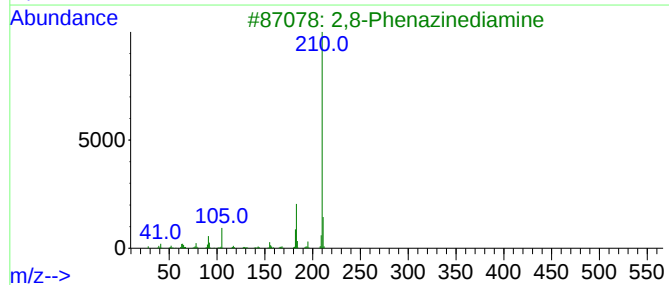
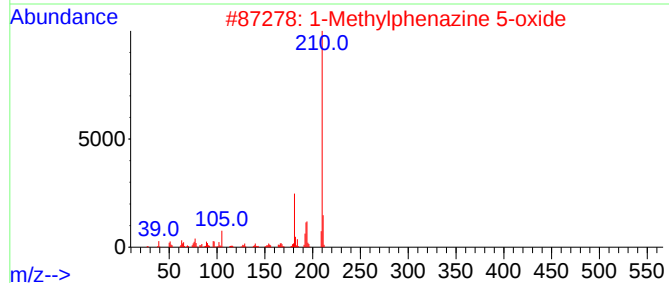
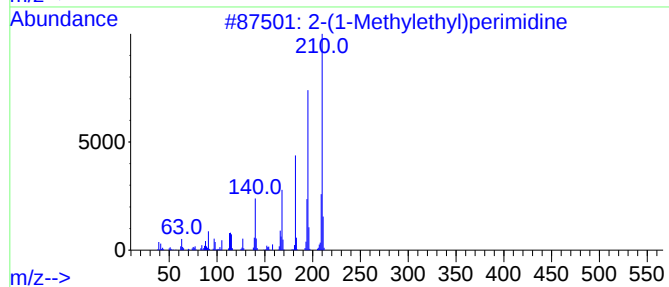
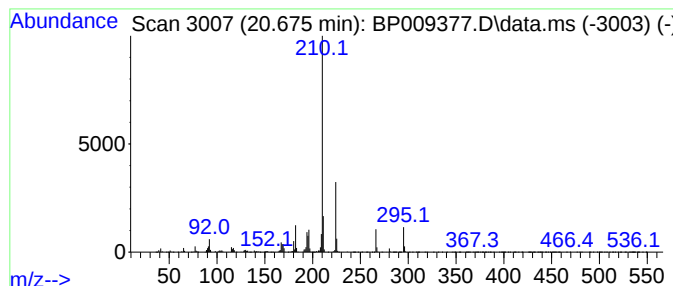
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 2-(1-Methylethyl)perimidine Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.675	23.67 ng	4902510	Chrysene-d12	21.333

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-(1-Methylethyl)perimidine	210	C14H14N2	028478-14-0	50
2		1-Methylphenazine 5-oxide	210	C13H10N2O	014202-95-0	50
3		2,8-Phenazinediamine	210	C12H10N4	007704-40-7	49
4		1,1'-Biphenyl, 3,3',4,4'-tetrame...	210	C16H18	004920-95-0	49
5		Anthracene, 1,2,3,4-tetrahydro-9...	210	C16H18	094573-50-9	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031122\
Data File : BP009377.D
Acq On : 11 Mar 2022 22:01
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Sample : N1886-01 20X
Misc :
ALS Vial : 20 Sample Multiplier: 1

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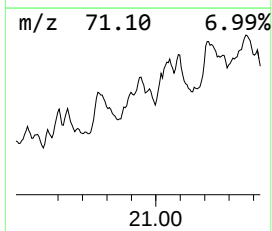
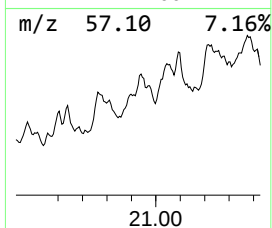
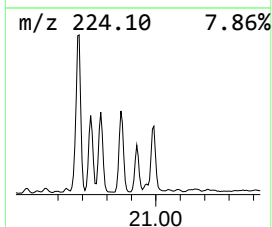
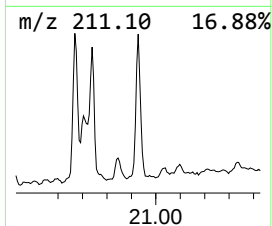
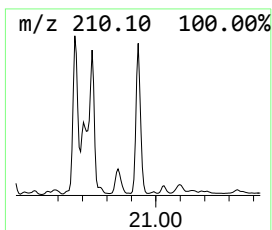
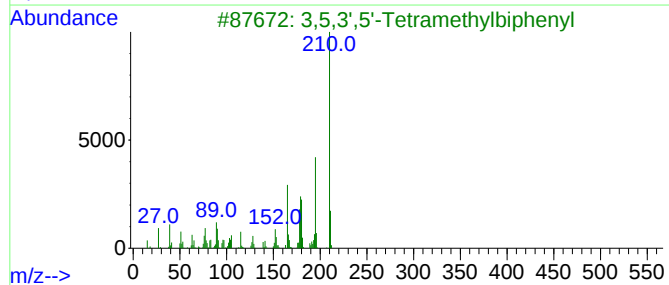
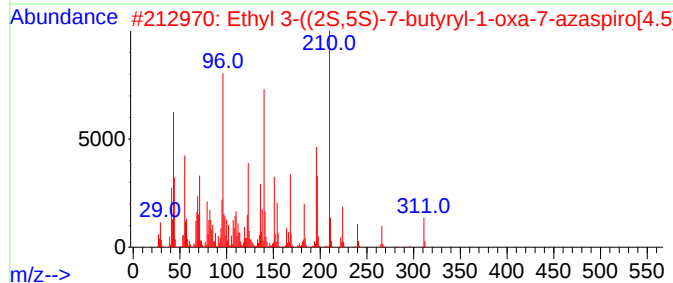
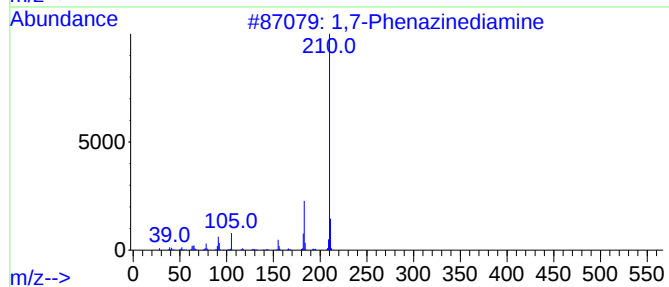
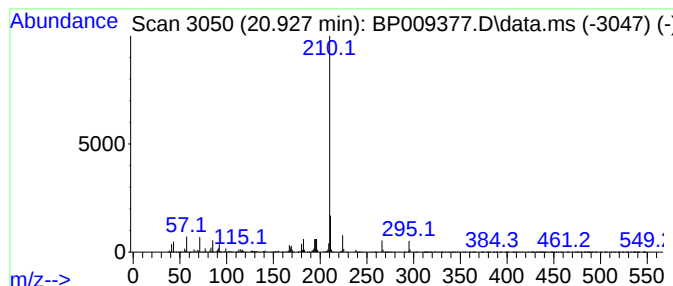
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1,7-Phenazinediamine Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.927	18.02 ng	3731820	Chrysene-d12	21.333

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,7-Phenazinediamine	210	C12H10N4	028124-29-0	64
2			Ethyl 3-((2S,5S)-7-butyryl-1-oxa...	311	C17H29NO4	1000483-96-5	64
3			3,5,3',5'-Tetramethylbiphenyl	210	C16H18	025570-02-9	64
4			1-Hydroxy-6-methylphenazine	210	C13H10N2O	014031-08-4	59
5			Carbamic acid, N-(2-butyl)-N-hex...	239	C14H25NO2	1000491-11-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031122\
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Acq On : 11 Mar 2022 22:01
Operator : CG/JU
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Misc :
ALS Vial : 20 Sample Multiplier: 1

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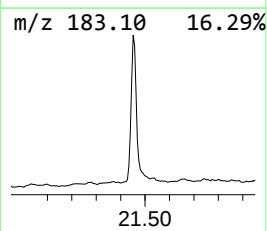
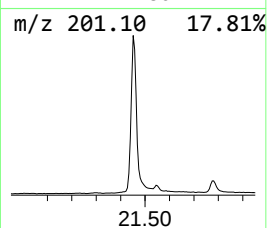
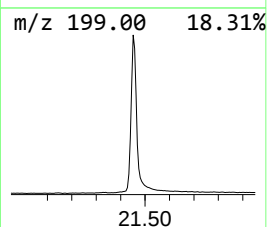
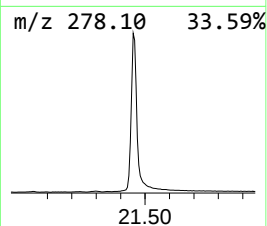
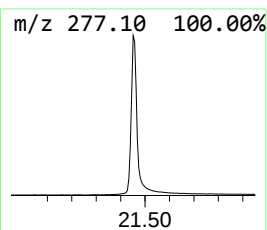
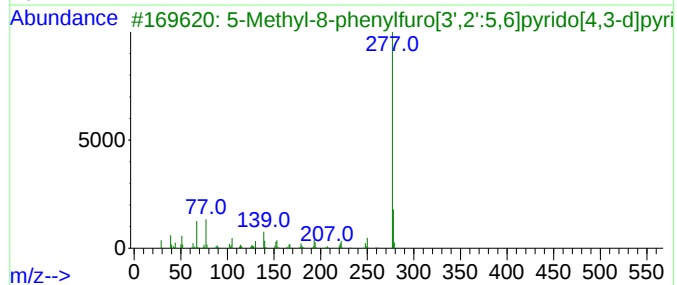
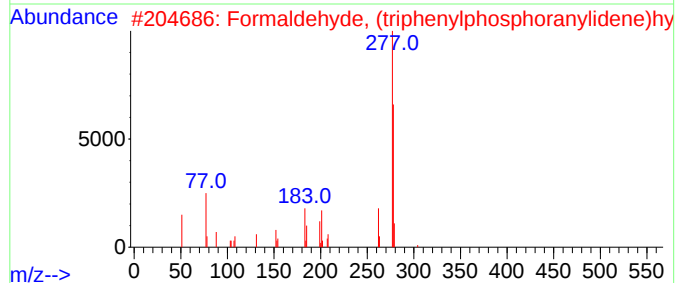
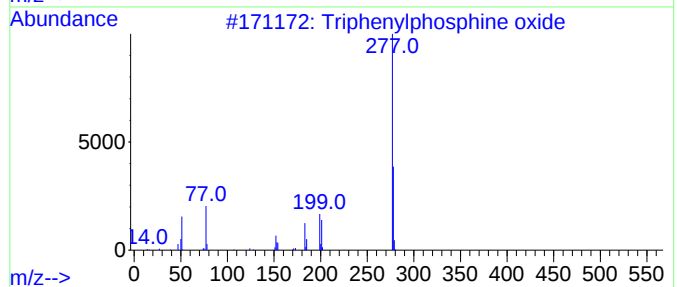
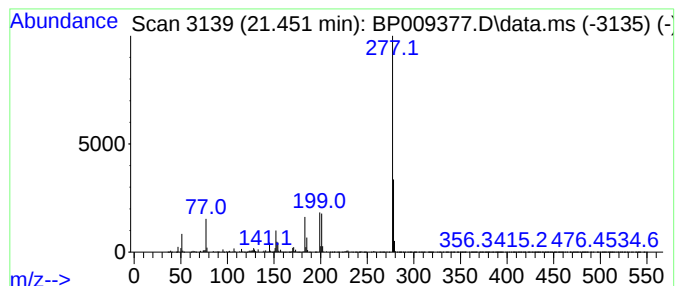
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Triphenylphosphine oxide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.451	70.06 ng	14512700	Chrysene-d12	21.333

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	94
2			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	47
3			5-Methyl-8-phenylfuro[3',2':5,6]...	277	C16H11N3O2	1000460-07-9	38
4			(E)-7-Benzylidene-1-azabicyclo[3...	293	C15H19N3S	1000483-83-4	35
5			2(1H)-Pyrimidinone, 5-methoxy-4,...	278	C17H14N2O2	028567-84-2	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031122\
Data File : BP009377.D
Acq On : 11 Mar 2022 22:01
Operator : CG/JU
Sample : N1886-01 20X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
2822-D

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Hexanol	3.893	5.2	ng	559026	1	7.816	2163660	20.0
Undecane	9.058	9.3	ng	1010520	1	7.816	2163660	20.0
unknown11.534	11.534	2.5	ng	426540	2	10.610	3453550	20.0
n-Hexadecanoic ...	18.128	3.5	ng	833929	4	17.222	4701930	20.0
Eicosane, 10-me...	18.416	2.3	ng	550542	4	17.222	4701930	20.0
Nonadecane	18.863	2.1	ng	503384	4	17.222	4701930	20.0
Heneicosane	19.028	2.2	ng	517347	4	17.222	4701930	20.0
unknown19.551	19.551	63.6	ng	13182300	5	21.333	4142760	20.0
Tridecane, 7-he...	19.592	2.6	ng	547320	5	21.333	4142760	20.0
2-(1-Methylethy...	20.675	23.7	ng	4902510	5	21.333	4142760	20.0
1,7-Phenazinedi...	20.927	18.0	ng	3731820	5	21.333	4142760	20.0
Triphenylphosph...	21.451	70.1	ng	14512700	5	21.333	4142760	20.0