

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102218\
 Data File : BN003251.D
 Acq On : 22 Oct 2018 22:08
 Operator : MJ/SJ
 Sample : J5568-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AK4

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 1 % 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_N\METHODS\SOM-EPA-BN101918MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.817	645	651	652	rBV	73906	105582	2.32%	0.450%
2	6.952	670	674	685	rVB	28649	51179	1.12%	0.218%
3	7.152	700	708	721	rBV	51573	89389	1.96%	0.381%
4	7.605	778	785	795	rBV	262686	443987	9.74%	1.892%
5	8.340	900	910	929	rVB	79780	170443	3.74%	0.726%
6	8.769	978	983	995	rBV	26846	56609	1.24%	0.241%
7	9.487	1100	1105	1114	rBV4	25653	51930	1.14%	0.221%
8	10.052	1194	1201	1215	rBV2	34266	95311	2.09%	0.406%
9	10.369	1249	1255	1267	rBV	424042	754368	16.55%	3.215%
10	10.546	1279	1285	1298	rBV2	38720	94191	2.07%	0.401%
11	13.663	1808	1815	1819	rBV	93956	135957	2.98%	0.579%
12	13.710	1819	1823	1833	rVB	72413	115228	2.53%	0.491%
13	13.940	1856	1862	1872	rBV	72471	114205	2.51%	0.487%
14	14.246	1907	1914	1926	rBV2	735669	1144689	25.12%	4.879%
15	14.569	1961	1969	1988	rBV4	22003	95054	2.09%	0.405%
16	15.251	2079	2085	2093	rBV	102360	163260	3.58%	0.696%
17	15.422	2109	2114	2125	rBV	45749	87748	1.93%	0.374%
18	16.416	2277	2283	2290	rBV2	51283	74260	1.63%	0.316%
19	16.998	2376	2382	2391	rBV	1007222	1439591	31.59%	6.135%
20	17.104	2391	2400	2410	rVB2	120576	203906	4.47%	0.869%
21	17.787	2510	2516	2529	rBV2	61768	141344	3.10%	0.602%
22	17.910	2531	2537	2558	rBV	1167732	1813912	39.81%	7.731%
23	18.575	2646	2650	2659	rBV4	41030	71981	1.58%	0.307%
24	19.045	2727	2730	2731	rVV2	52195	58738	1.29%	0.250%
25	19.086	2731	2737	2751	rVV5	169371	514060	11.28%	2.191%
26	19.192	2751	2755	2768	rVB	235855	366524	8.04%	1.562%
27	19.410	2787	2792	2802	rVB	158241	246716	5.41%	1.051%
28	20.051	2897	2901	2905	rVV	68958	104523	2.29%	0.445%
29	20.175	2919	2922	2924	rBV4	37301	45798	1.01%	0.195%
30	20.204	2924	2927	2937	rVB6	51312	99713	2.19%	0.425%
31	20.957	3051	3055	3057	rBV3	62487	91686	2.01%	0.391%
32	21.210	3093	3098	3109	rVB	1443421	1925260	42.25%	8.205%
33	21.357	3119	3123	3127	rVB	92232	106846	2.34%	0.455%
34	21.780	3191	3195	3200	rVB	188137	227570	4.99%	0.970%

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Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	22.233	3268	3272	3278	rVB	321674	403109	8.85%	1.718%
36	22.357	3289	3293	3299	rVB	224172	282208	6.19%	1.203%
37	22.727	3351	3356	3361	rVB	364093	530360	11.64%	2.260%
38	23.274	3443	3449	3459	rVB	538294	910788	19.99%	3.882%
39	23.416	3466	3473	3489	rBV	965148	2109970	46.30%	8.992%
40	23.892	3549	3554	3562	rVB	313599	587897	12.90%	2.506%
41	24.363	3627	3634	3649	rVB	109711	444022	9.74%	1.892%
42	24.633	3670	3680	3703	rBV5	1130700	4556845	100.00%	19.421%
43	24.974	3730	3738	3765	rBV3	482375	2230100	48.94%	9.504%
44	25.439	3813	3817	3819	rBV2	63977	106873	2.35%	0.455%

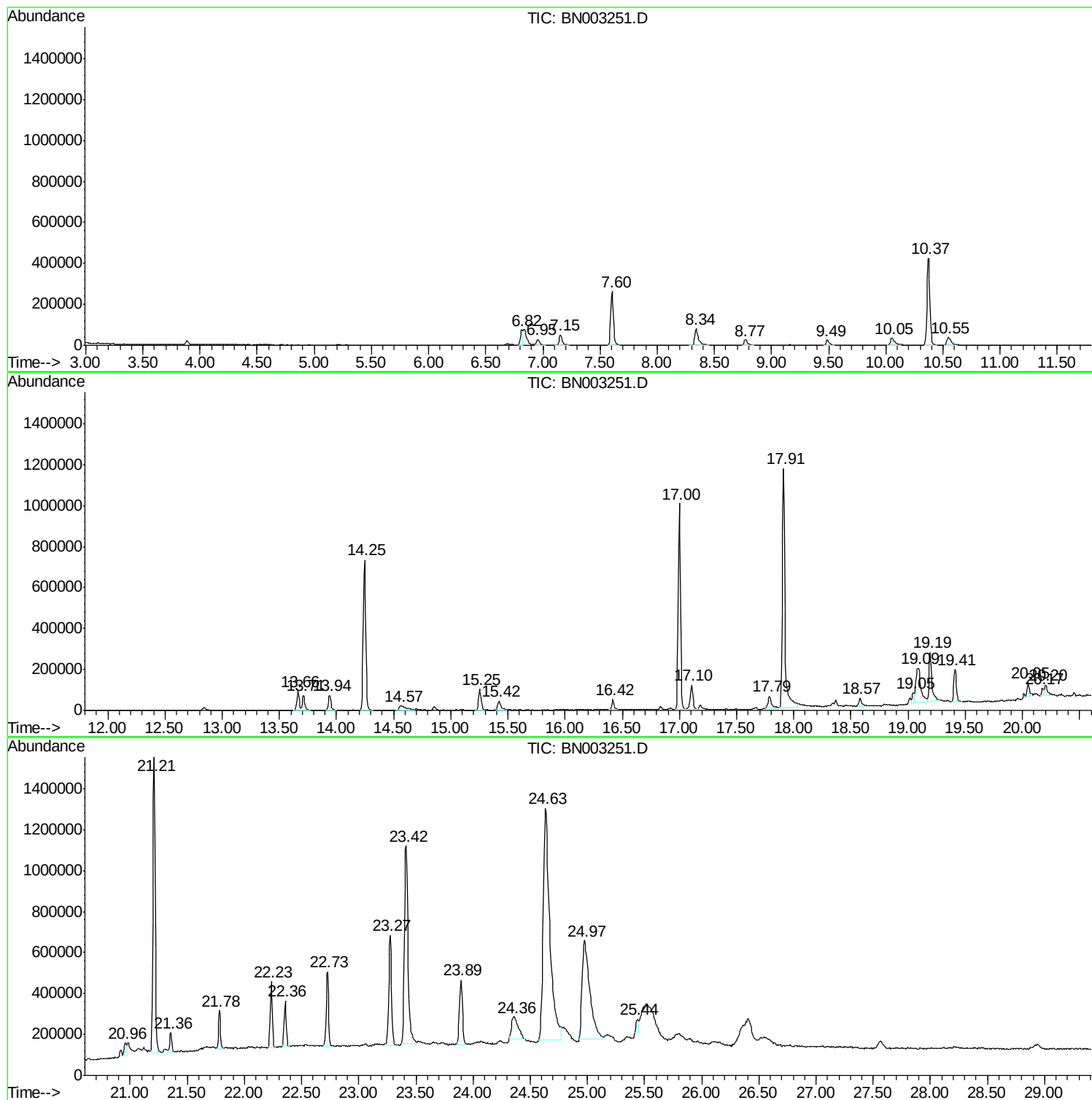
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN101918MA.M
Quant Title : SVOA CALIBRATION

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



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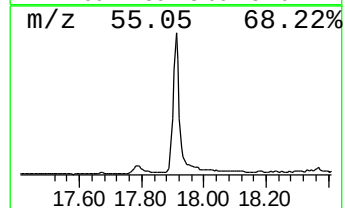
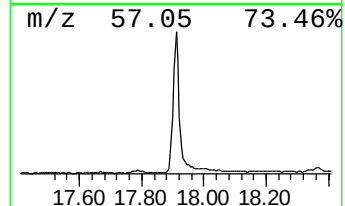
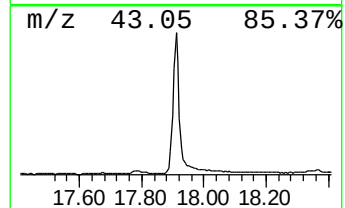
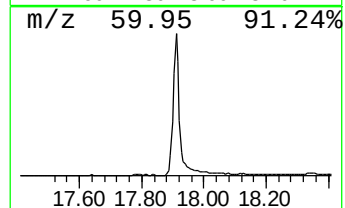
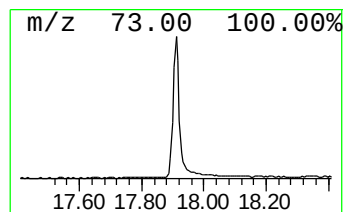
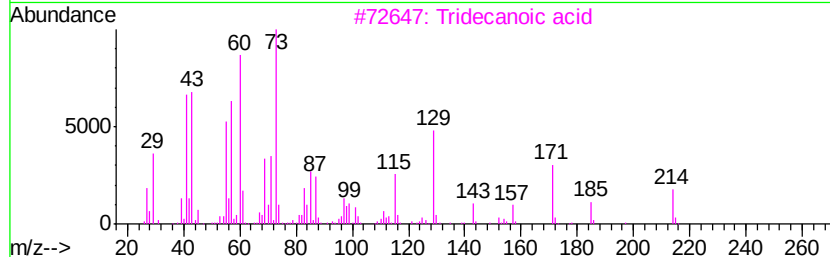
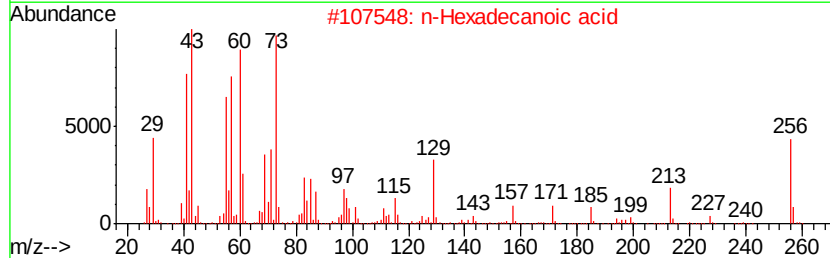
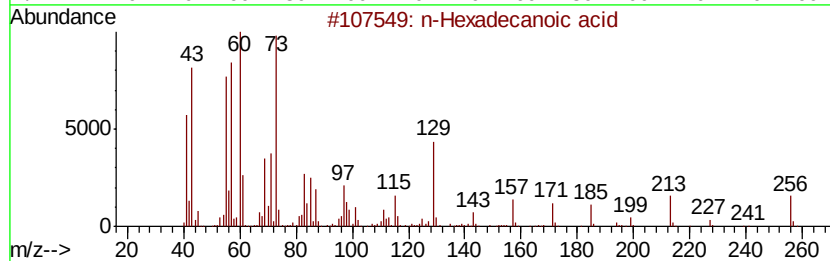
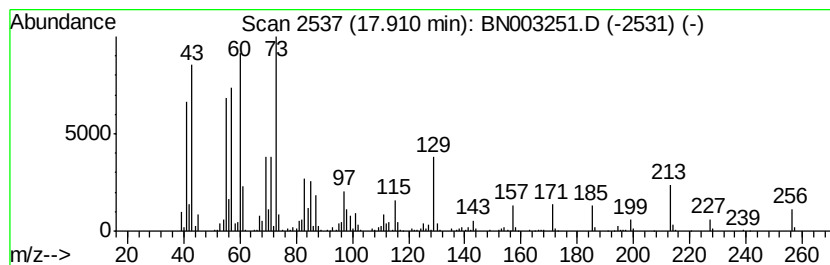
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN101918MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.91	25.20 ng/ul	1813910	Phenanthrene-d10	17.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		Tridecanoic acid	214	C13H26O2	000638-53-9	93
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
5		Tridecanoic acid	214	C13H26O2	000638-53-9	86



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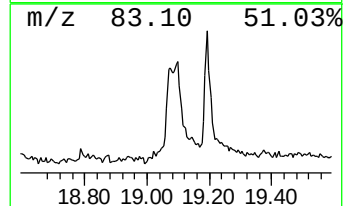
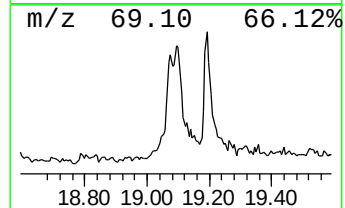
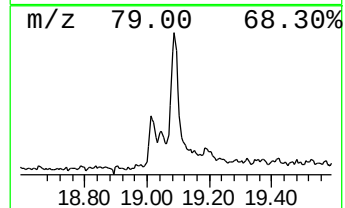
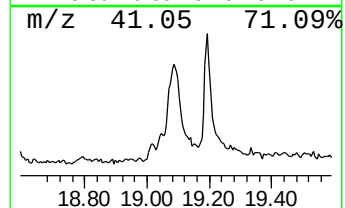
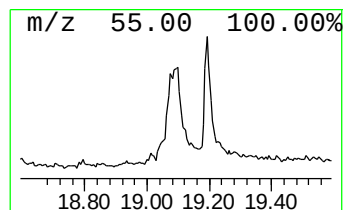
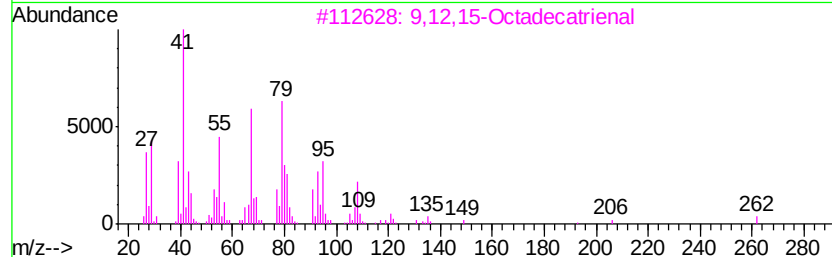
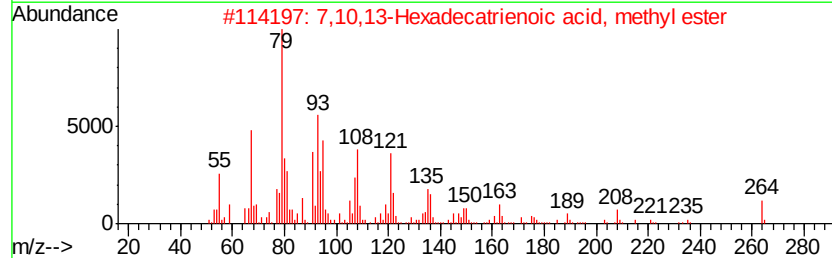
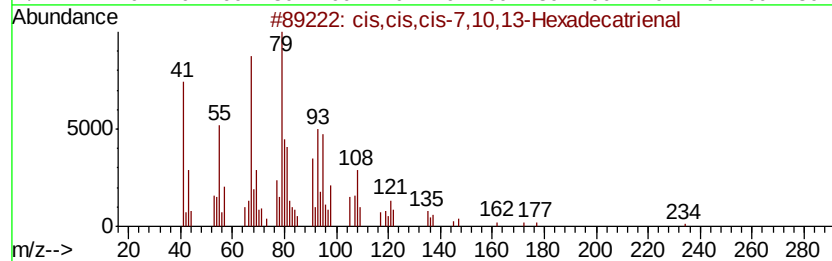
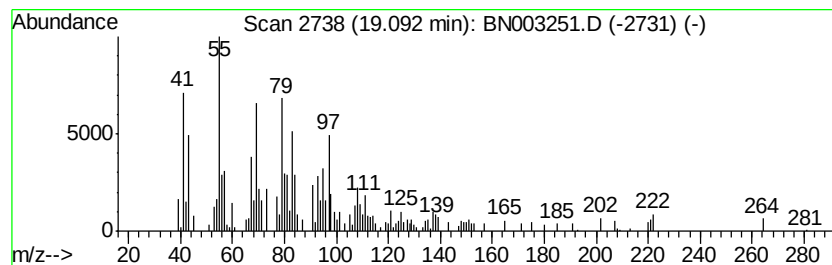
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Peak Number 2 cis,cis,cis-7,10,13-Hexadec... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.09	7.14 ng/ul	514060	Phenanthrene-d10	17.00		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis,cis,cis-7,10,13-Hexadecatrienal	234	C16H26O	056797-43-4	53	
2	7,10,13-Hexadecatrienoic acid, m...	264	C17H28O2	056554-30-4	49	
3	9,12,15-Octadecatrienal	262	C18H30O	026537-71-3	49	
4	9,12,15-Octadecatrienoic acid, (...	278	C18H30O2	000463-40-1	45	
5	9,12,15-Octadecatrien-1-ol, (Z,Z...	264	C18H32O	000506-44-5	43	



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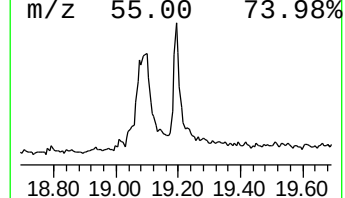
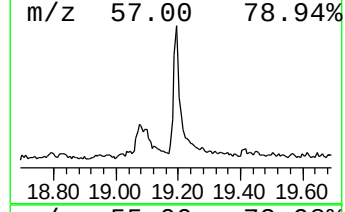
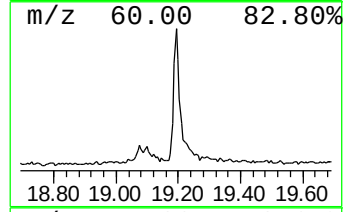
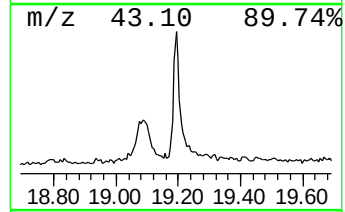
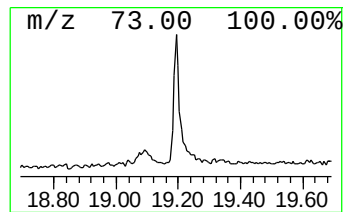
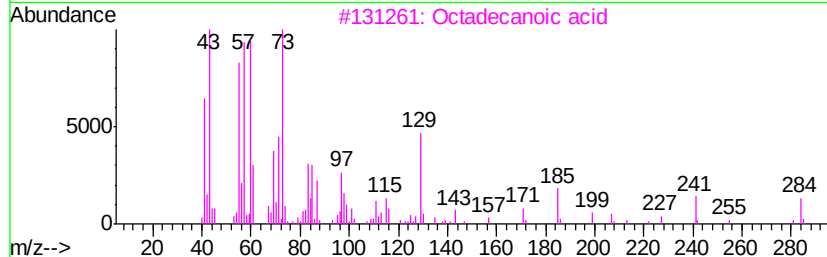
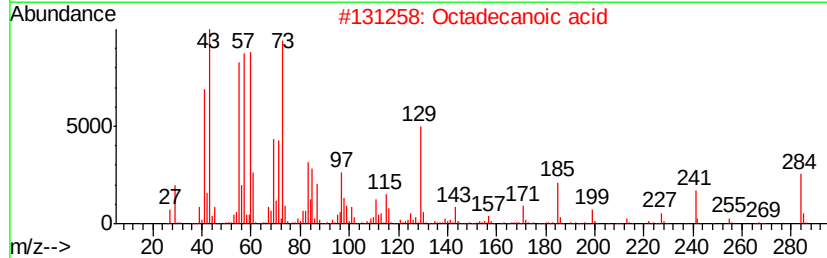
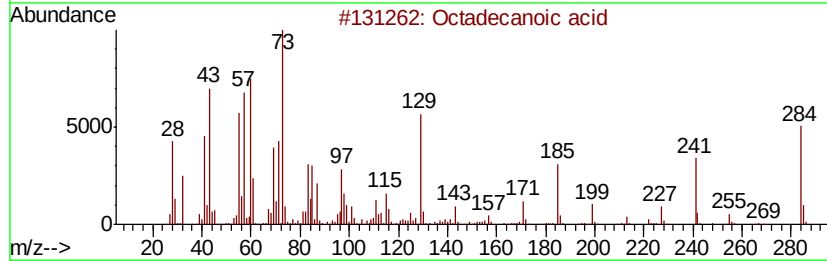
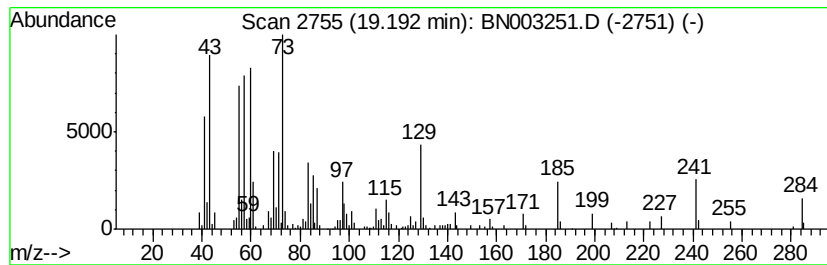
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.19	3.81 ng/ul	366524	Chrysene-d12	21.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid	284	C18H36O2	000057-11-4	99
3		Octadecanoic acid	284	C18H36O2	000057-11-4	96
4		Octadecanoic acid	284	C18H36O2	000057-11-4	95
5		Octadecanoic acid	284	C18H36O2	000057-11-4	93



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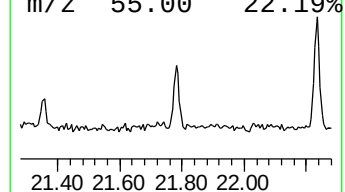
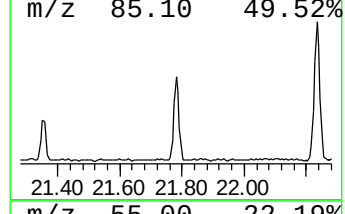
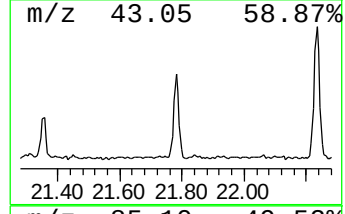
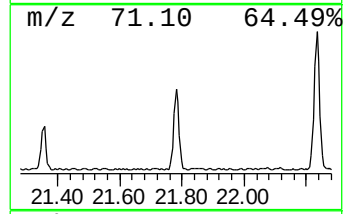
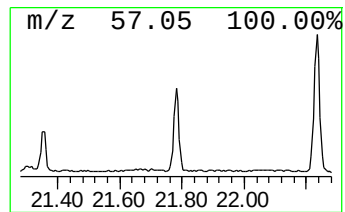
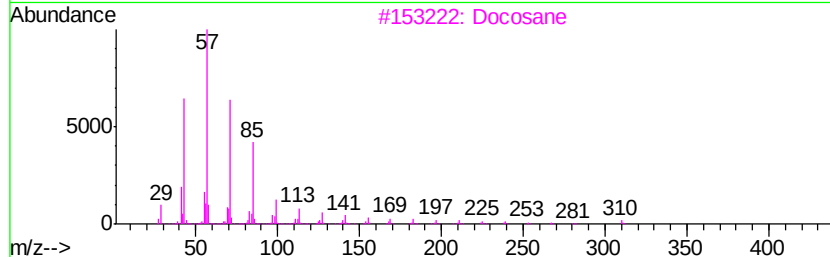
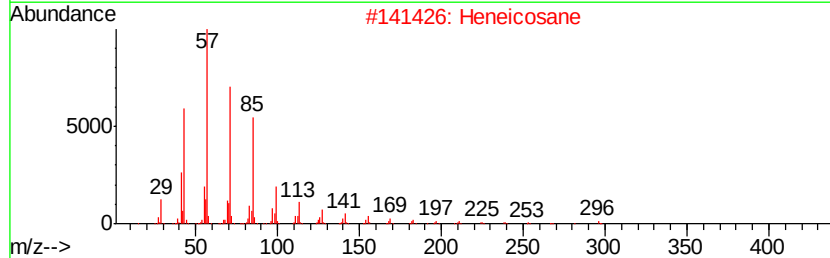
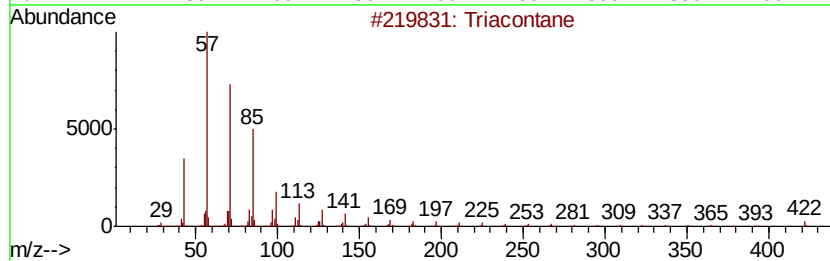
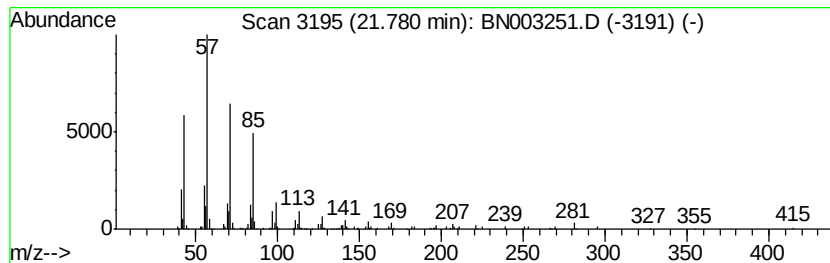
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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.78	2.36 ng/ul	227570	Chrysene-d12	21.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triacontane	422	C30H62	000638-68-6	90
2		Heneicosane	296	C21H44	000629-94-7	90
3		Docosane	310	C22H46	000629-97-0	90
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	87
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



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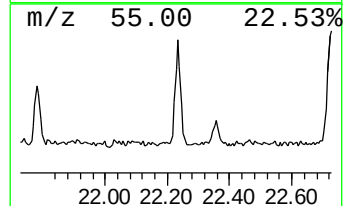
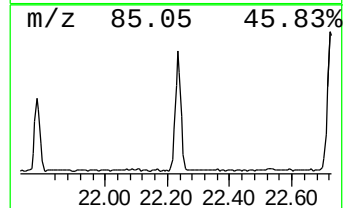
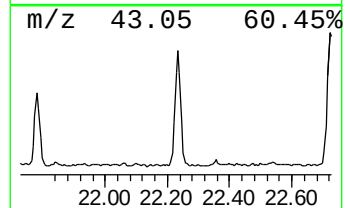
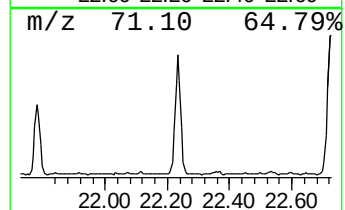
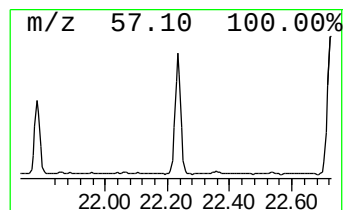
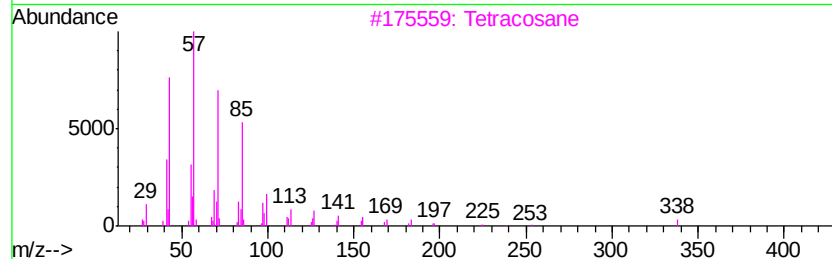
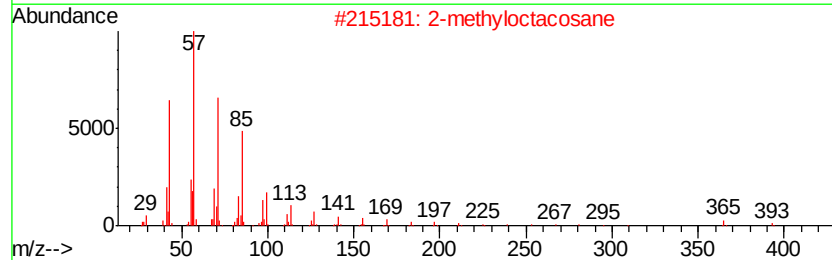
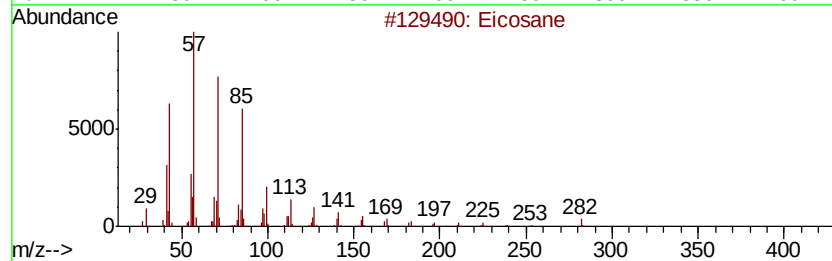
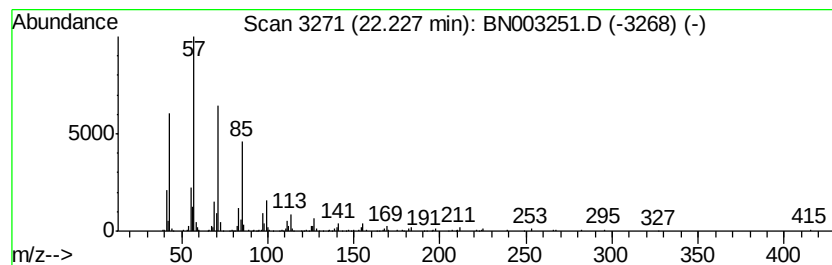
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN101918MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.23	4.19 ng/ul	403109	Chrysene-d12	21.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	96
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Octacosane	394	C28H58	000630-02-4	91



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Acq On : 22 Oct 2018 22:08
Operator : MJ/SJ
Sample : J5568-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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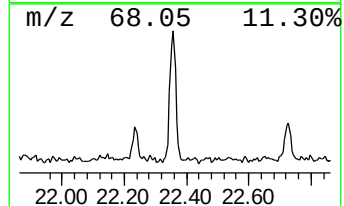
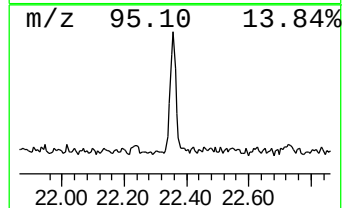
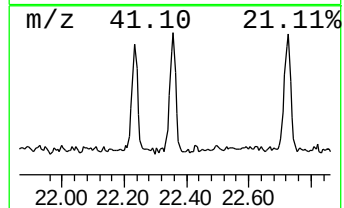
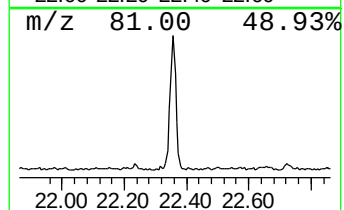
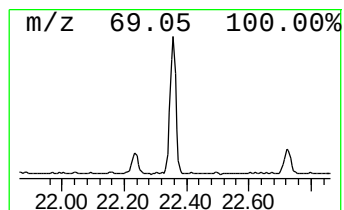
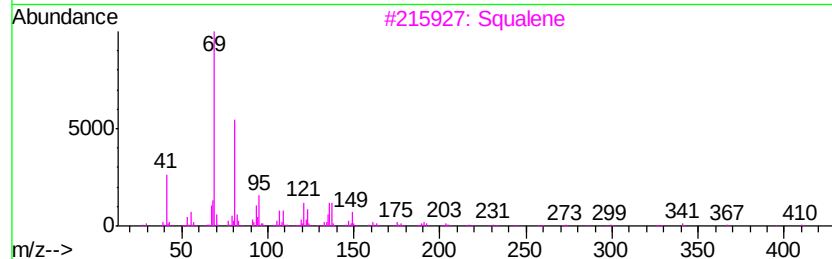
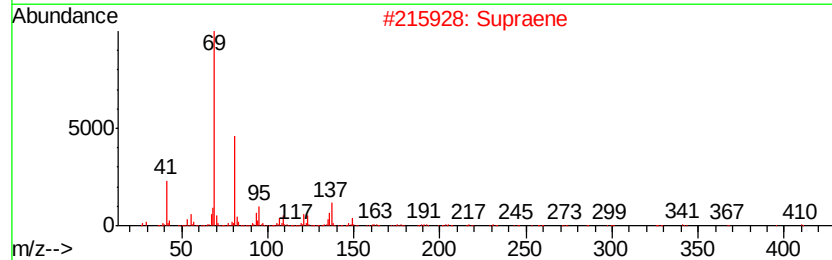
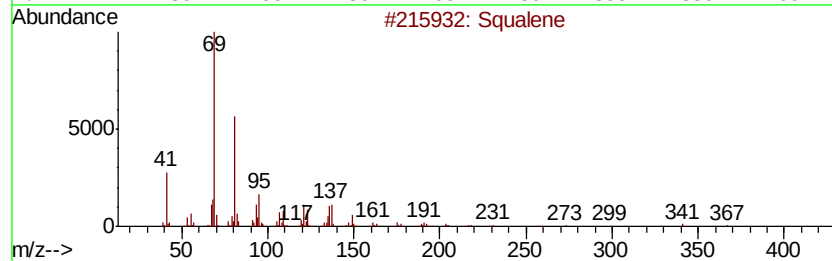
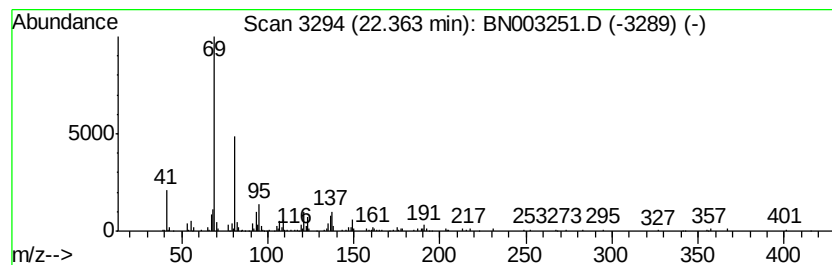
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Squalene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.36	2.68 ng/ul	282208	Perylene-d12	23.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	91
2		Supraene	410	C30H50	007683-64-9	91
3		Squalene	410	C30H50	000111-02-4	91
4		Squalene	410	C30H50	000111-02-4	91
5		Squalene	410	C30H50	000111-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102218\
Data File : BN003251.D
Acq On : 22 Oct 2018 22:08
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Sample : J5568-09
Misc :
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Instrument :
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ClientSampleId :
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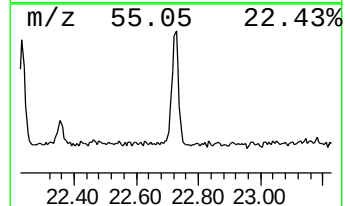
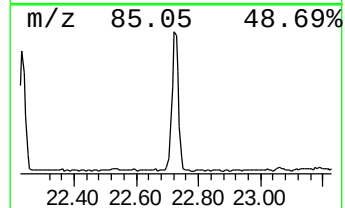
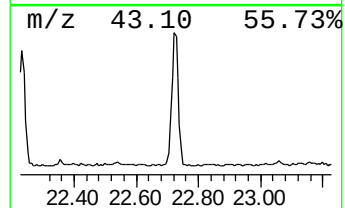
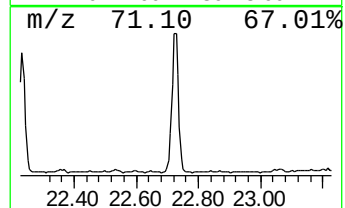
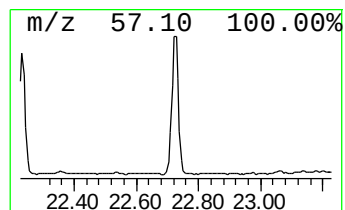
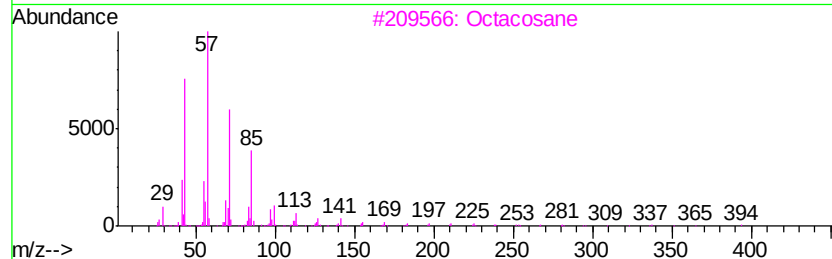
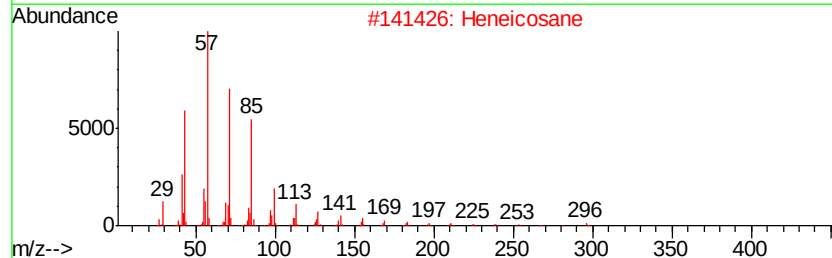
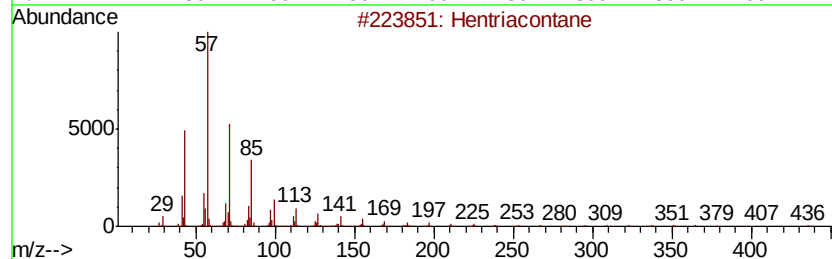
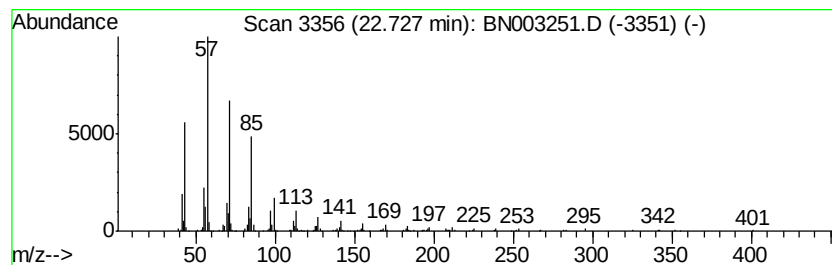
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.73	5.03 ng/ul	530360	Perylene-d12	23.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hentriacontane	437	C31H64	000630-04-6	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Octacosane	394	C28H58	000630-02-4	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102218\
Data File : BN003251.D
Acq On : 22 Oct 2018 22:08
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Sample : J5568-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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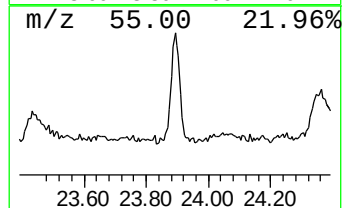
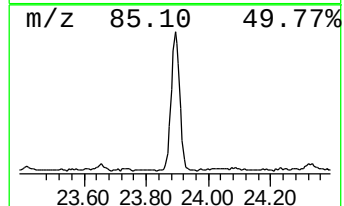
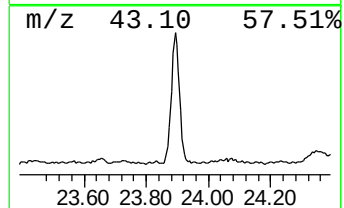
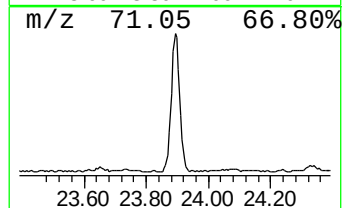
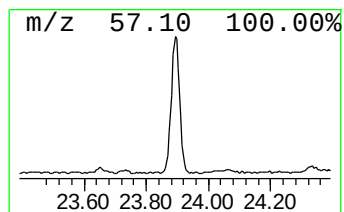
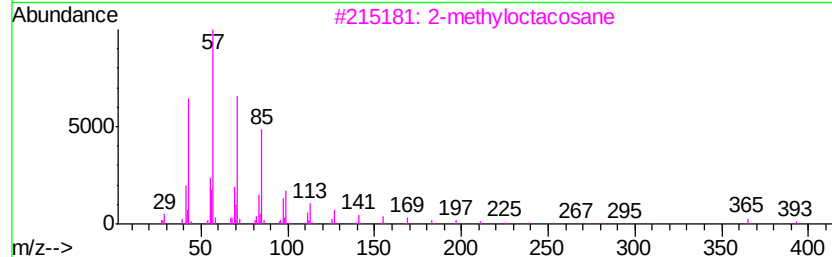
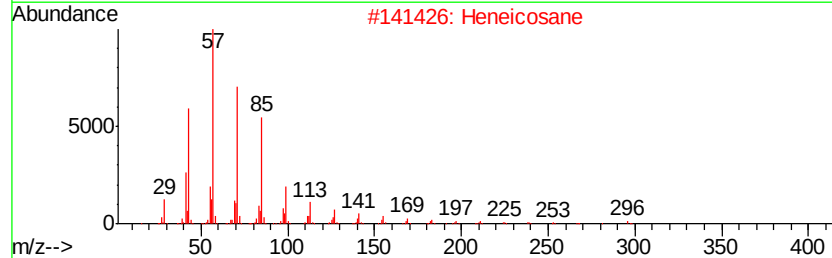
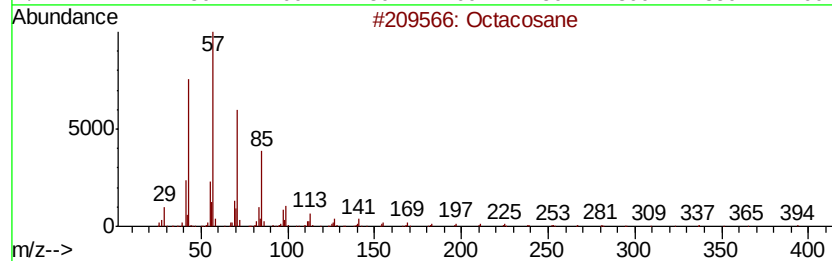
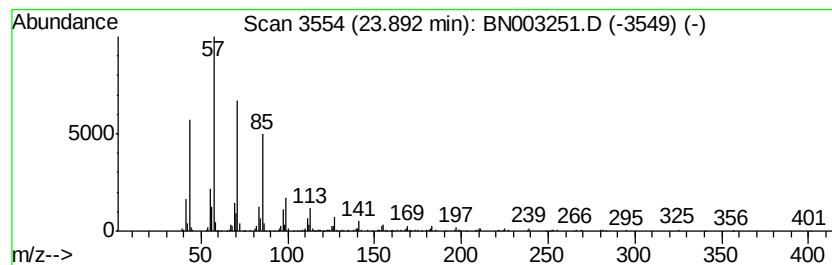
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.89	5.57 ng/ul	587897	Perylene-d12	23.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
5		Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102218\
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Sample : J5568-09
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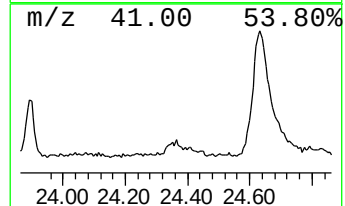
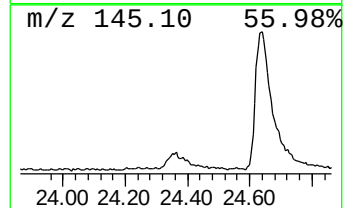
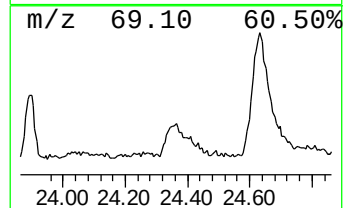
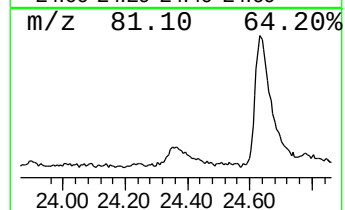
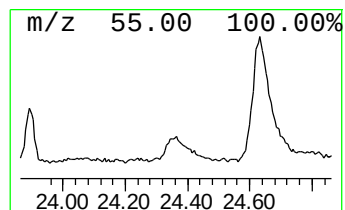
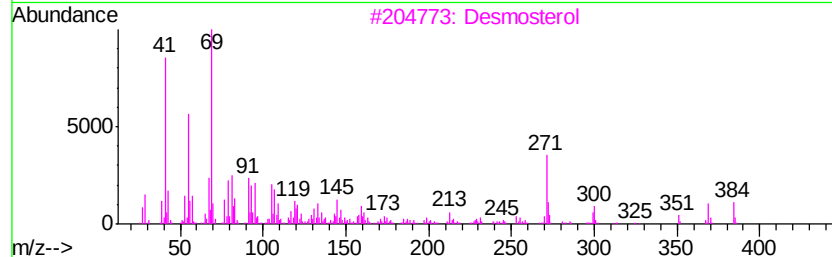
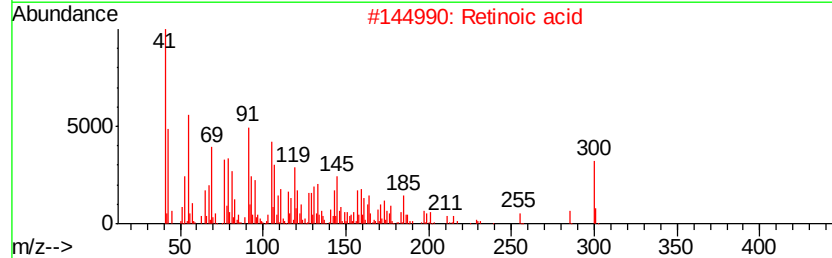
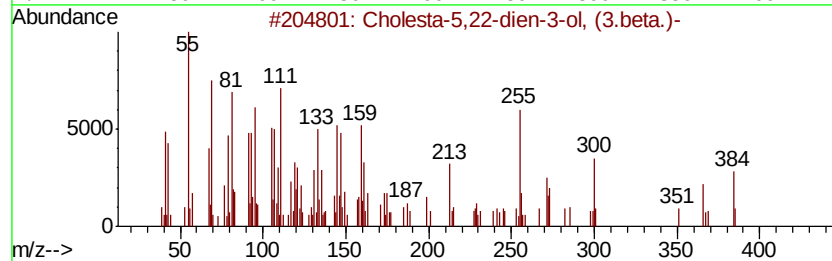
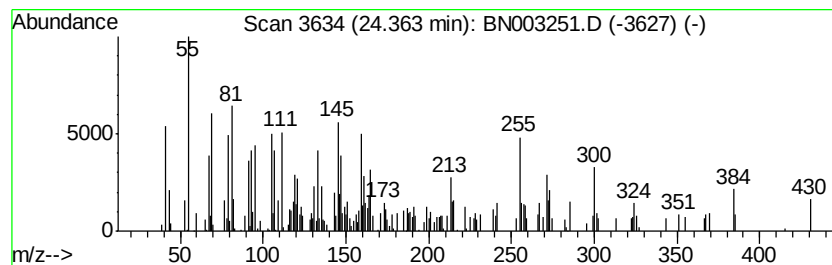
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Cholesta-5,22-dien-3-ol, (3... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
24.36	4.21 ng/ul	444022	Perylene-d12	23.41		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cholesta-5,22-dien-3-ol, (3.beta.)-	384	C27H44O	092218-20-7	90	
2	Retinoic acid	300	C20H28O2	000302-79-4	25	
3	Desmosterol	384	C27H44O	000313-04-2	25	
4	1,4,2,5-Dioxadiazine, 3,6-bis(2,...	322	C20H22N2O2	020434-87-1	11	
5	1,3,4-Thiadiazol-2-amine, 5-trif...	273	C11H10F3N3S	284488-18-2	10	



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102218\
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Sample : J5568-09
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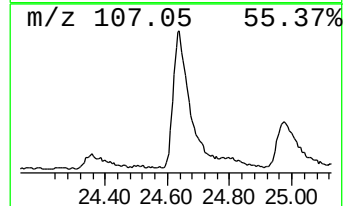
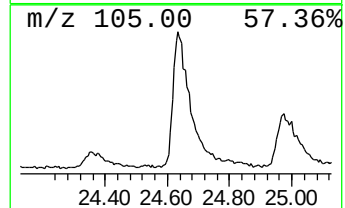
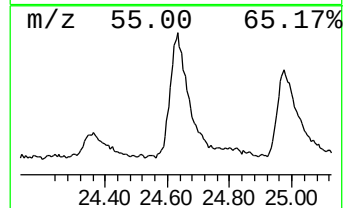
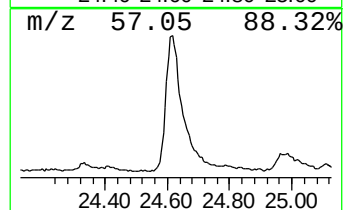
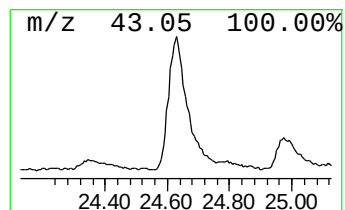
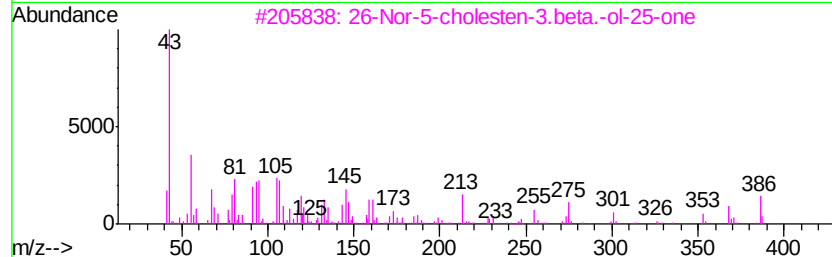
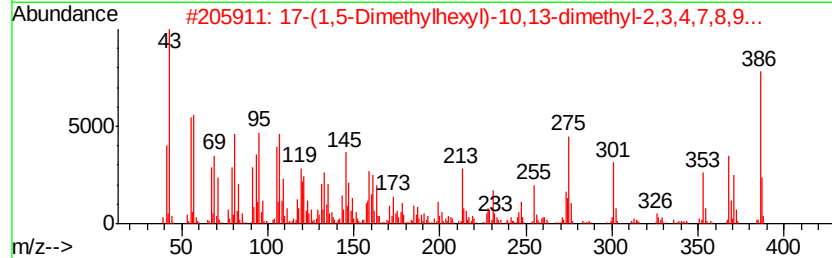
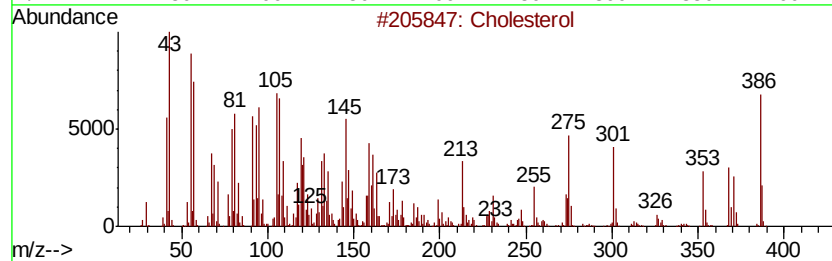
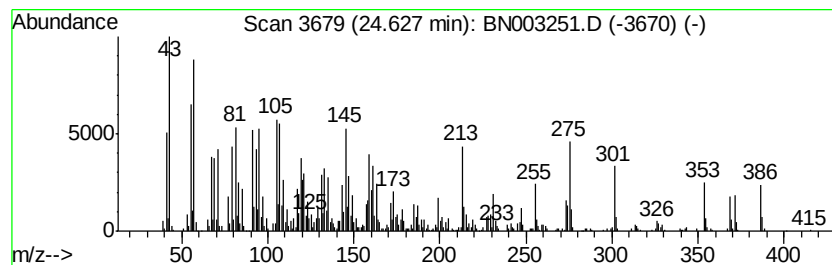
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Cholesterol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.63	43.19 ng/ul	4556850	Perylene-d12	23.41
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Cholesterol		386 C27H46O	000057-88-5 99
2	17-(1,5-Dimethylhexyl)-10,13-dim...		386 C27H46O	1000210-38-4 99
3	26-Nor-5-cholesten-3.beta.-ol-25...		386 C26H42O2	007494-34-0 99
4	Cholesterol		386 C27H46O	000057-88-5 95
5	26-Nor-5-cholesten-3.beta.-ol-25...		386 C26H42O2	007494-34-0 70



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102218\
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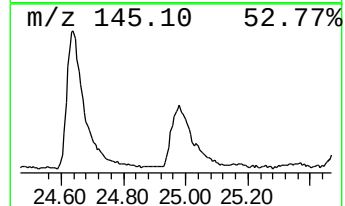
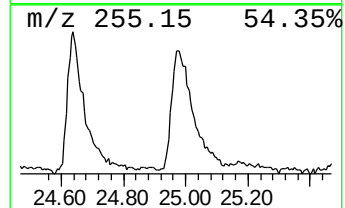
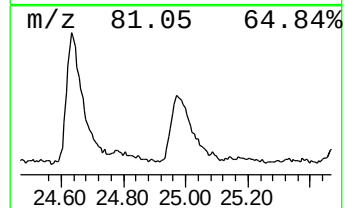
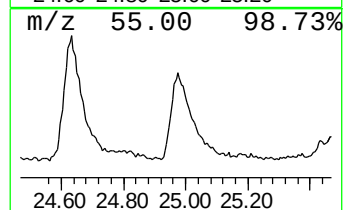
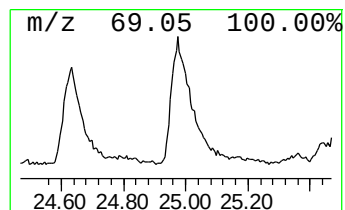
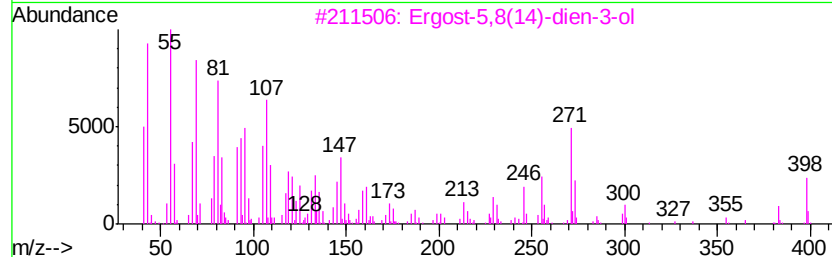
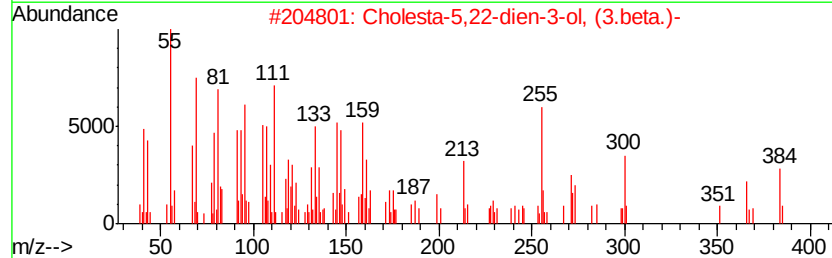
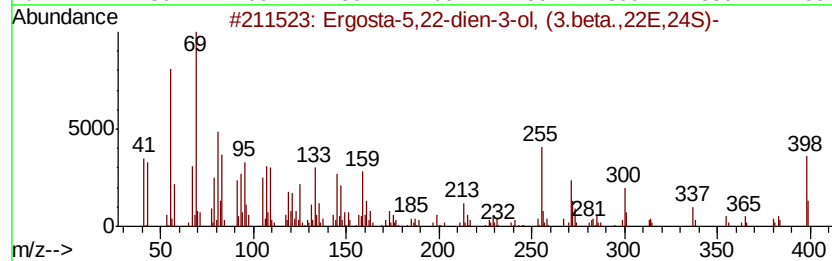
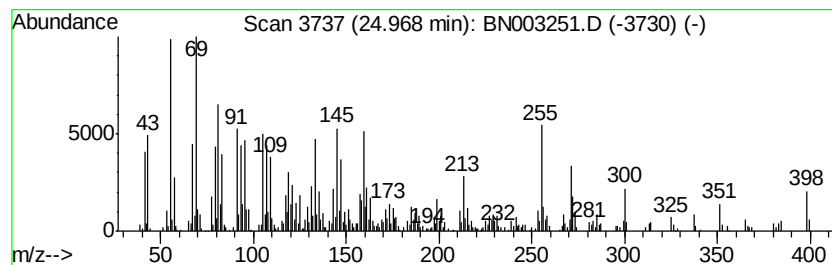
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Ergosta-5,22-dien-3-ol, (3.... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.97	21.14 ng/ul	2230100	Perylene-d12	23.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ergosta-5,22-dien-3-ol, (3.beta....	398	C28H46O	017472-78-5	96
2		Cholesta-5,22-dien-3-ol, (3.beta.)-	384	C27H44O	092218-20-7	70
3		Ergost-5,8(14)-dien-3-ol	398	C28H46O	177962-83-3	64
4		Ergosta-5,7-dien-3-ol, (3.beta.)-	398	C28H46O	000516-79-0	64
5		Ergosta-5,22-dien-3-ol, (3.beta....	398	C28H46O	000474-67-9	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN102218\
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 Sample : J5568-09
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 ALS Vial : 10 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN101918MA.M
 Quant Title : SVOA 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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cis,cis,cis-7,10,...	19.09	7.1	ng/ul	514060	4	17.00	1439590	20.0
Octadecanoic acid	19.19	3.8	ng/ul	366524	5	21.21	1925260	20.0
(DEL) Alkane: Str...	21.78	2.4	ng/ul	227570	5	21.21	1925260	20.0
(DEL) Alkane: Str...	22.23	4.2	ng/ul	403109	5	21.21	1925260	20.0
Squalene	22.36	2.7	ng/ul	282208	6	23.41	2109970	20.0
(DEL) Alkane: Str...	22.73	5.0	ng/ul	530360	6	23.41	2109970	20.0
(DEL) Alkane: Str...	23.89	5.6	ng/ul	587897	6	23.41	2109970	20.0
Cholesta-5,22-die...	24.36	4.2	ng/ul	444022	6	23.41	2109970	20.0
Cholesterol	24.63	43.2	ng/ul	4556850	6	23.41	2109970	20.0
Ergosta-5,22-dien...	24.97	21.1	ng/ul	2230100	6	23.41	2109970	20.0