

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100924\
 Data File : BM047960.D
 Acq On : 10 Oct 2024 01:17
 Operator : RC/JU
 Sample : P4187-09
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EY070

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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_M\Methods\SFAM-EPA-BM092724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM047960.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.240	40	43	49	rVB	52373	66314	1.12%	0.128%
2	3.663	112	115	128	rBV	145355	243109	4.12%	0.469%
3	3.987	166	170	173	rVB3	11905	19112	0.32%	0.037%
4	4.440	242	247	254	rBV2	19486	47177	0.80%	0.091%
5	4.781	302	305	310	rVB5	11245	16202	0.27%	0.031%
6	4.905	322	326	331	rVB2	15550	24601	0.42%	0.047%
7	6.051	515	521	529	rVB3	22585	35754	0.61%	0.069%
8	6.346	566	571	572	rBV5	6878	9275	0.16%	0.018%
9	6.957	667	675	687	rBV	241846	441573	7.47%	0.852%
10	7.116	696	702	714	rBV	690560	1110247	18.79%	2.141%
11	7.322	730	737	748	rBV	826297	1321314	22.37%	2.548%
12	7.422	750	754	758	rVB7	3792	6564	0.11%	0.013%
13	7.787	809	816	825	rVV	800614	1288596	21.81%	2.485%
14	7.857	825	828	836	rVV6	11132	25665	0.43%	0.049%
15	8.157	873	879	882	rBV8	7723	10757	0.18%	0.021%
16	8.281	897	900	903	rVB5	5542	7524	0.13%	0.015%
17	8.492	930	936	953	rBV	629832	1094212	18.52%	2.110%
18	8.710	968	973	976	rBV6	5244	9420	0.16%	0.018%
19	8.940	1005	1012	1028	rBV	765120	1329122	22.50%	2.563%
20	9.069	1029	1034	1039	rVB2	33058	49665	0.84%	0.096%
21	9.375	1082	1086	1091	rBV	16127	24581	0.42%	0.047%
22	9.516	1106	1110	1113	rBV7	5896	10746	0.18%	0.021%
23	9.663	1126	1135	1146	rBV	633126	1082108	18.32%	2.087%
24	10.204	1219	1227	1245	rBV	963437	1709560	28.94%	3.297%
25	10.581	1283	1291	1302	rBV	1058892	1757214	29.74%	3.389%
26	10.716	1306	1314	1339	rBV	734680	1440404	24.38%	2.778%
27	10.975	1356	1358	1366	rVB8	6445	13423	0.23%	0.026%
28	11.875	1506	1511	1515	rBV3	8367	11486	0.19%	0.022%
29	12.175	1558	1562	1567	rBV3	11515	20245	0.34%	0.039%
30	13.157	1726	1729	1735	rVB6	7459	9927	0.17%	0.019%
31	13.251	1742	1745	1751	rVB7	6836	10814	0.18%	0.021%
32	13.827	1837	1843	1853	rBV	1689292	2451799	41.50%	4.729%
33	14.116	1882	1892	1903	rBV	2048355	3009643	50.94%	5.805%
34	14.422	1937	1944	1952	rBV	1528701	2192365	37.11%	4.228%
35	14.627	1972	1979	1994	rBV	138631	306208	5.18%	0.591%
36	14.845	2012	2016	2021	rVB7	7671	11392	0.19%	0.022%

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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % 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_M\Methods\SFAM-EPA-BM092724.MA.M
 Title : SVOA CALIBRATION

37	15.227	2076	2081	2085	rBV7	4601	8124	0.14%	0.016%
38	15.416	2105	2113	2126	rBV	2644721	3829792	64.83%	7.386%
39	15.527	2126	2132	2144	rBV	818634	1175777	19.90%	2.268%
40	15.657	2149	2154	2159	rVB3	12668	20993	0.36%	0.040%
41	15.774	2169	2174	2181	rBV	226684	319526	5.41%	0.616%
42	15.986	2206	2210	2212	rVB4	6244	7679	0.13%	0.015%
43	16.198	2243	2246	2252	rVB6	4969	6783	0.11%	0.013%
44	16.339	2263	2270	2280	rBV	41075	64013	1.08%	0.123%
45	17.163	2403	2410	2418	rBV	1763661	2555717	43.26%	4.929%
46	17.257	2420	2426	2441	rVV2	3089398	4358772	73.78%	8.407%
47	17.451	2454	2459	2462	rVB3	6144	7614	0.13%	0.015%
48	17.563	2474	2478	2480	rBV5	4771	6932	0.12%	0.013%
49	17.827	2521	2523	2528	rVB5	5618	7631	0.13%	0.015%
50	18.051	2556	2561	2569	rBV2	74173	119958	2.03%	0.231%
51	18.357	2610	2613	2616	rBV5	12252	18530	0.31%	0.036%
52	18.904	2701	2706	2713	rBV6	19616	34504	0.58%	0.067%
53	19.115	2738	2742	2747	rBV3	36920	52872	0.89%	0.102%
54	19.180	2750	2753	2759	rBV2	37826	53145	0.90%	0.102%
55	19.327	2775	2778	2786	rBV7	40661	65529	1.11%	0.126%
56	19.545	2809	2815	2828	rBV	3982185	5486305	92.86%	10.581%
57	20.539	2981	2984	2989	rVB	99244	106490	1.80%	0.205%
58	21.380	3122	3127	3140	rBV	1918380	3008282	50.92%	5.802%
59	24.168	3592	3601	3613	rBV	2492734	5907855	100.00%	11.394%
60	24.374	3627	3636	3648	rVB	1333819	3409056	57.70%	6.575%

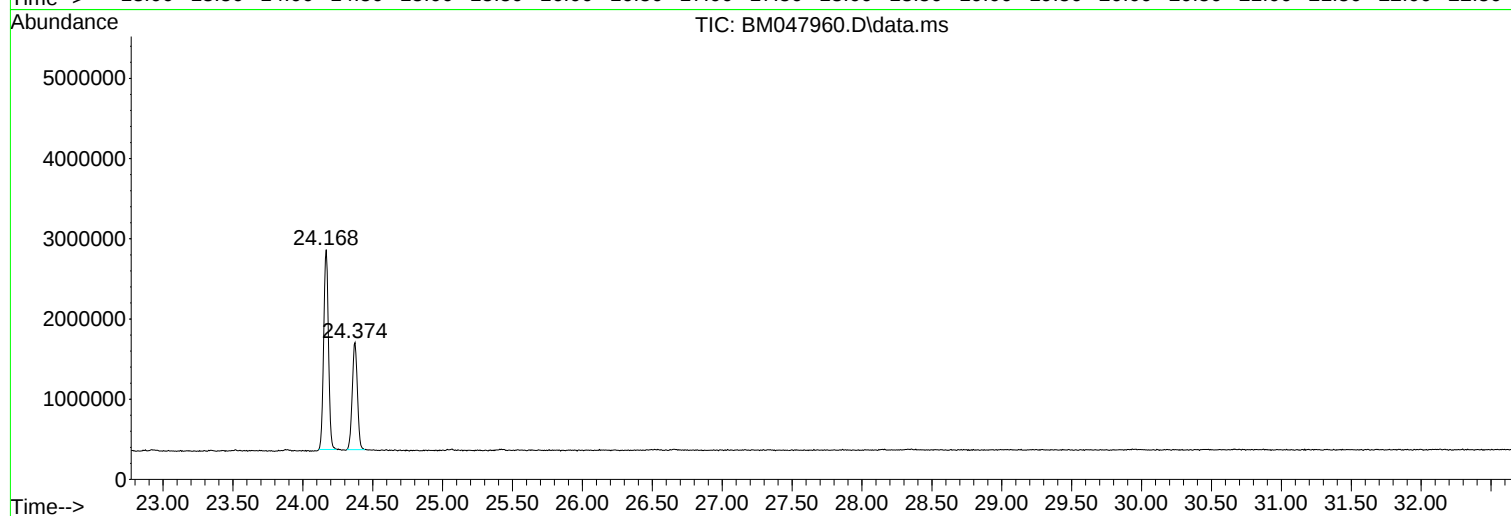
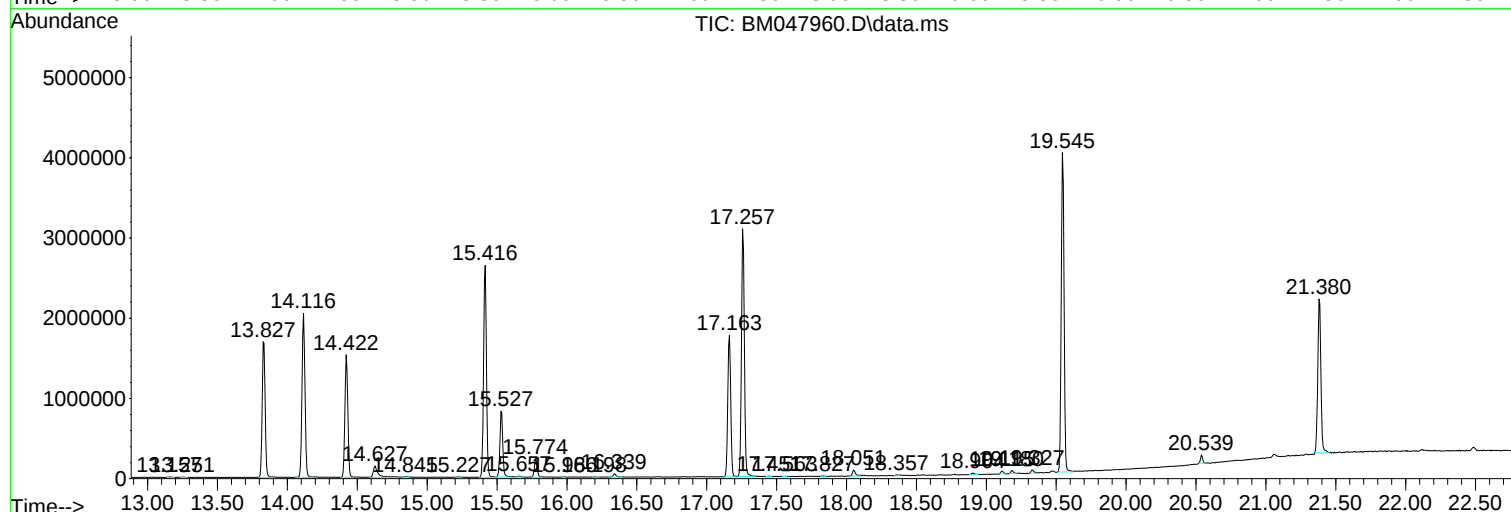
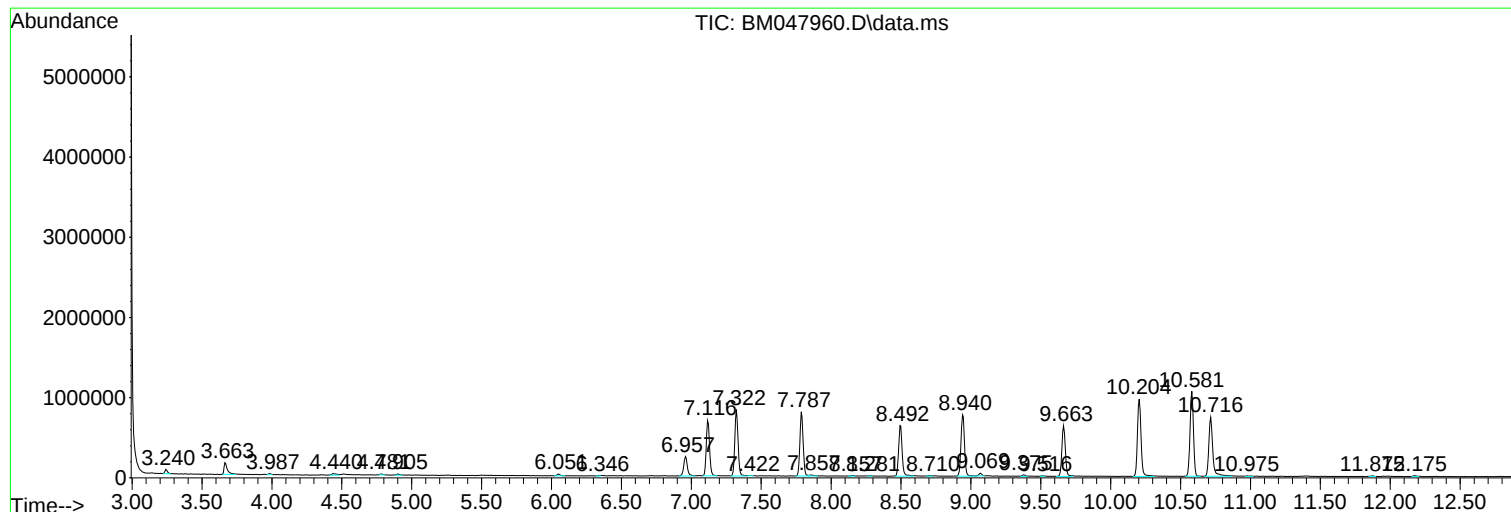
Sum of corrected areas: 51849997

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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



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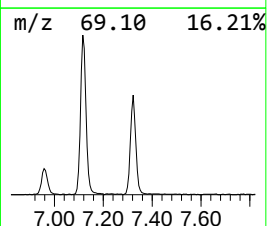
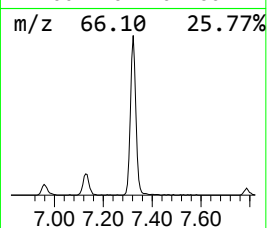
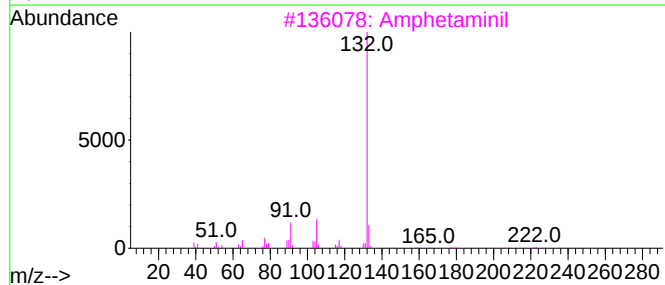
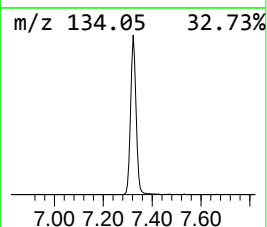
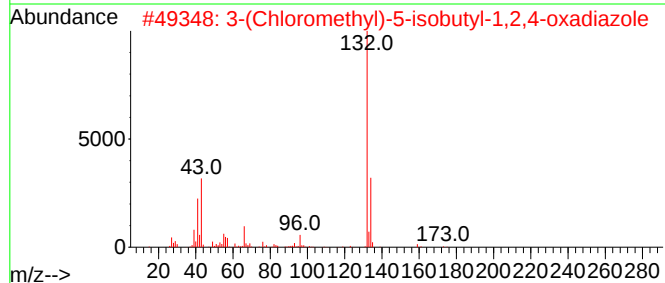
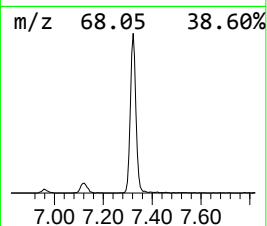
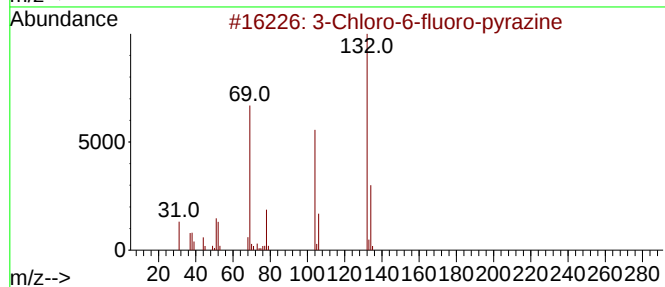
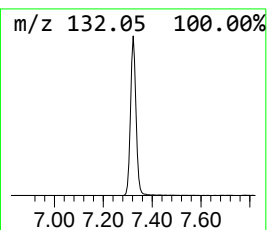
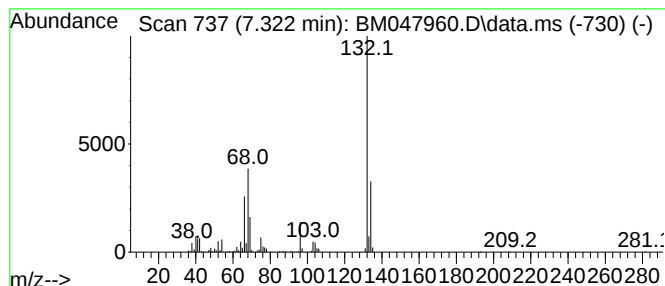
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.322	20.51 ng/ul	1321310	1,4-Dichlorobenzene-d4	7.787

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			3-(Chloromethyl)-5-isobutyl-1,2,...	174	C7H11ClN2O	189130-85-6	25
3			Amphetaminil	250	C17H18N2	017590-01-1	10
4			Pyrazino[2,3-b]pyrazine	132	C6H4N4	000255-53-8	9
5			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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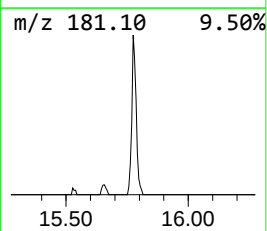
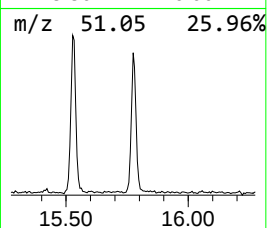
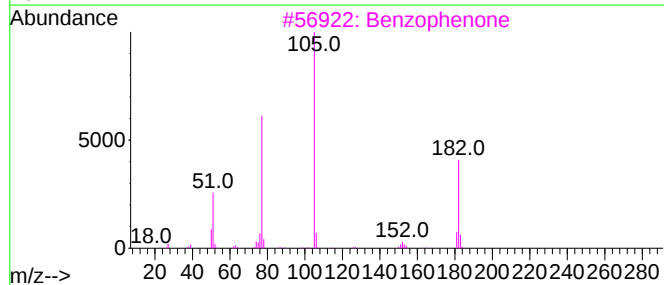
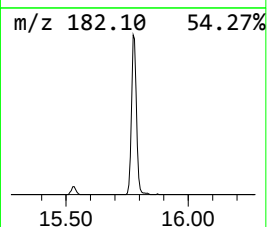
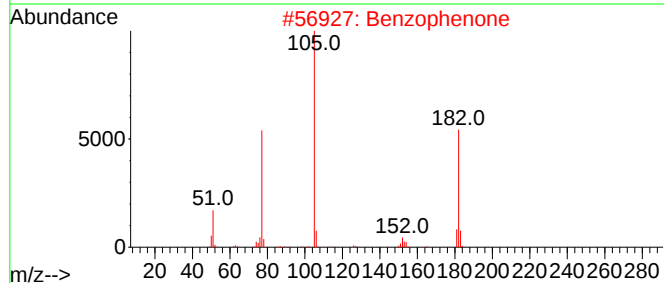
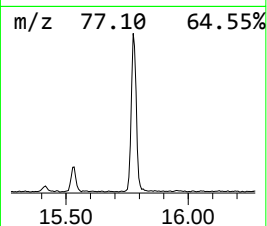
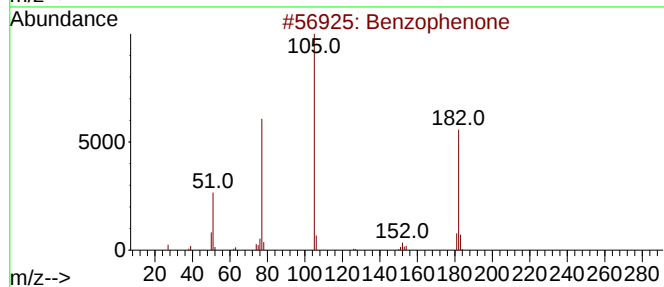
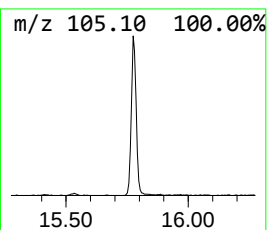
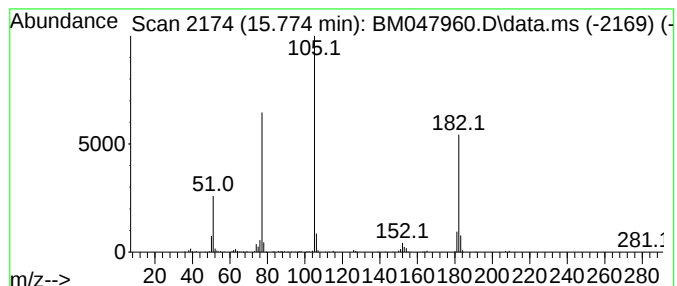
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.774	2.91 ng/ul	319526	Acenaphthene-d10	14.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	93
5			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49



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Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	7.322	20.5	ng/u1	1321310	1	7.787	1288600	20.0
Benzophenone	15.774	2.9	ng/u1	319526	3	14.422	2192370	20.0