

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013258.D
 Acq On : 22 Dec 2022 23:00
 Operator : CG/JU
 Sample : N6015-19
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXZ8

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_P\Methods\SFAM-EPA-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013258.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	21	25	33	rVB	253999	319594	0.61%	0.112%
2	3.758	94	97	112	rBV	2284994	3488834	6.64%	1.222%
3	3.869	113	116	120	rVB2	43877	60172	0.11%	0.021%
4	4.087	150	153	160	rVB	50283	70704	0.13%	0.025%
5	5.040	309	315	321	rVB2	32367	58553	0.11%	0.021%
6	5.258	347	352	362	rBV	116573	196825	0.37%	0.069%
7	6.616	575	583	591	rVB	477828	709045	1.35%	0.248%
8	7.110	659	667	679	rVB	3839990	6082174	11.57%	2.130%
9	7.275	686	695	707	rVB	3186023	4881875	9.29%	1.709%
10	7.375	707	712	717	rBV	82936	123524	0.23%	0.043%
11	7.475	722	729	741	rBV	4007415	6250024	11.89%	2.188%
12	7.946	800	809	816	rBV	2361656	3653491	6.95%	1.279%
13	8.657	918	930	939	rBV	3632182	5903934	11.23%	2.067%
14	9.116	1000	1008	1016	rBV	3889345	6225124	11.84%	2.180%
15	9.516	1070	1076	1083	rBV2	66345	141538	0.27%	0.050%
16	9.704	1099	1108	1114	rBV	371807	580670	1.10%	0.203%
17	9.840	1122	1131	1146	rBV	3275561	5417801	10.31%	1.897%
18	10.069	1164	1170	1178	rVB3	118405	236438	0.45%	0.083%
19	10.381	1215	1223	1238	rBV	4986284	8178792	15.56%	2.864%
20	10.763	1279	1288	1295	rBV	3102903	5197487	9.89%	1.820%
21	10.898	1301	1311	1327	rBV	2628507	4710974	8.96%	1.649%
22	11.069	1335	1340	1344	rVB3	47693	71413	0.14%	0.025%
23	12.116	1513	1518	1523	rBV4	62642	91652	0.17%	0.032%
24	12.340	1551	1556	1561	rBV2	80760	124459	0.24%	0.044%
25	13.345	1721	1727	1730	rBV5	39211	66212	0.13%	0.023%
26	13.404	1733	1737	1742	rVV	74419	117077	0.22%	0.041%
27	13.457	1742	1746	1751	rVV6	39422	70709	0.13%	0.025%
28	13.510	1751	1755	1759	rVV2	40574	66510	0.13%	0.023%
29	13.910	1817	1823	1830	rBV	206165	307613	0.59%	0.108%
30	13.998	1830	1838	1845	rVV	7993038	11044599	21.01%	3.867%
31	14.063	1846	1849	1852	rVV4	38762	53489	0.10%	0.019%
32	14.110	1853	1857	1861	rVB6	44618	64242	0.12%	0.022%
33	14.286	1879	1887	1895	rBV	9150783	13848407	26.35%	4.849%
34	14.475	1916	1919	1925	rVB	50189	72180	0.14%	0.025%
35	14.545	1925	1931	1932	rBV5	34760	52759	0.10%	0.018%
36	14.586	1932	1938	1945	rVB	4371053	6482298	12.33%	2.270%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Title : SVOA CALIBRATION

37	14.651	1945	1949	1953	rBV3	122329	178135	0.34%	0.062%
38	14.786	1966	1972	1977	rBV	2537290	3919346	7.46%	1.372%
39	14.834	1977	1980	1988	rVB2	343702	529633	1.01%	0.185%
40	14.939	1994	1998	2003	rVV2	114957	153266	0.29%	0.054%
41	14.992	2003	2007	2012	rVV4	56490	94976	0.18%	0.033%
42	15.045	2012	2016	2020	rVB5	49587	67829	0.13%	0.024%
43	15.498	2085	2093	2098	rBV2	170397	254358	0.48%	0.089%
44	15.586	2100	2108	2121	rBV	11115428	16488219	31.37%	5.773%
45	15.704	2121	2128	2136	rBV	3102541	4119751	7.84%	1.442%
46	15.828	2144	2149	2154	rVB2	95349	152945	0.29%	0.054%
47	15.945	2160	2169	2176	rVB	975228	1352180	2.57%	0.473%
48	16.028	2178	2183	2184	rBV	107442	123414	0.23%	0.043%
49	16.157	2199	2205	2212	rBV2	209253	307051	0.58%	0.108%
50	16.292	2224	2228	2235	rVB8	48270	98261	0.19%	0.034%
51	16.433	2249	2252	2257	rBV7	44136	75080	0.14%	0.026%
52	16.516	2263	2266	2271	rVB5	52306	73494	0.14%	0.026%
53	16.728	2298	2302	2308	rBV2	121307	196618	0.37%	0.069%
54	16.839	2317	2321	2325	rBV4	39315	61075	0.12%	0.021%
55	17.227	2380	2387	2394	rBV4	127441	239700	0.46%	0.084%
56	17.292	2395	2398	2401	rVV4	84567	102515	0.20%	0.036%
57	17.345	2401	2407	2417	rVV	4748268	7104494	13.52%	2.487%
58	17.445	2417	2424	2432	rVB2	10553454	14876505	28.30%	5.209%
59	17.557	2440	2443	2446	rVB3	53147	61231	0.12%	0.021%
60	17.698	2463	2467	2470	rBV3	102541	152621	0.29%	0.053%
61	17.963	2509	2512	2515	rVB3	47656	56902	0.11%	0.020%
62	17.998	2515	2518	2522	rVB4	47036	59059	0.11%	0.021%
63	18.104	2532	2536	2540	rBV4	77771	105087	0.20%	0.037%
64	18.157	2541	2545	2550	rBV	307384	453588	0.86%	0.159%
65	18.216	2551	2555	2565	rBV	2007506	2824411	5.37%	0.989%
66	18.504	2600	2604	2608	rBV5	86330	144959	0.28%	0.051%
67	18.551	2608	2612	2616	rVB2	93284	136475	0.26%	0.048%
68	18.698	2634	2637	2641	rVB3	75396	82239	0.16%	0.029%
69	18.827	2656	2659	2662	rBV4	63308	72090	0.14%	0.025%
70	19.198	2719	2722	2726	rBV	106184	148500	0.28%	0.052%
71	19.251	2728	2731	2736	rVB	141914	193117	0.37%	0.068%
72	19.351	2737	2748	2752	rBV3	477439	1383943	2.63%	0.485%
73	19.445	2760	2764	2770	rBV	413663	562030	1.07%	0.197%
74	19.674	2798	2803	2817	rVB	11745090	16156234	30.74%	5.657%
75	20.616	2960	2963	2967	rVB	210568	286877	0.55%	0.100%
76	21.127	3045	3050	3060	rVB2	1811727	2613053	4.97%	0.915%

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Title : SVOA CALIBRATION

77	21.433	3097	3102	3107	rBV	4826070	6511918	12.39%	2.280%
78	22.068	3207	3210	3222	rVB2	729352	1108699	2.11%	0.388%
79	23.127	3385	3390	3396	rBV2	1709819	2861840	5.44%	1.002%
80	23.757	3489	3497	3511	rVB	8552851	18130844	34.49%	6.348%
81	23.909	3517	3523	3531	rVB2	3319956	6840475	13.01%	2.395%
82	24.527	3622	3628	3637	rVB	3051463	6481216	12.33%	2.269%
83	24.627	3639	3645	3660	rVB	2566821	6769381	12.88%	2.370%
84	26.439	3947	3953	3974	rVB3	1544801	5535985	10.53%	1.938%
85	27.462	4119	4127	4148	rVB3	2098063	7722895	14.69%	2.704%
86	28.515	4292	4306	4329	rVB2	12866083	52563804	100.00%	18.404%

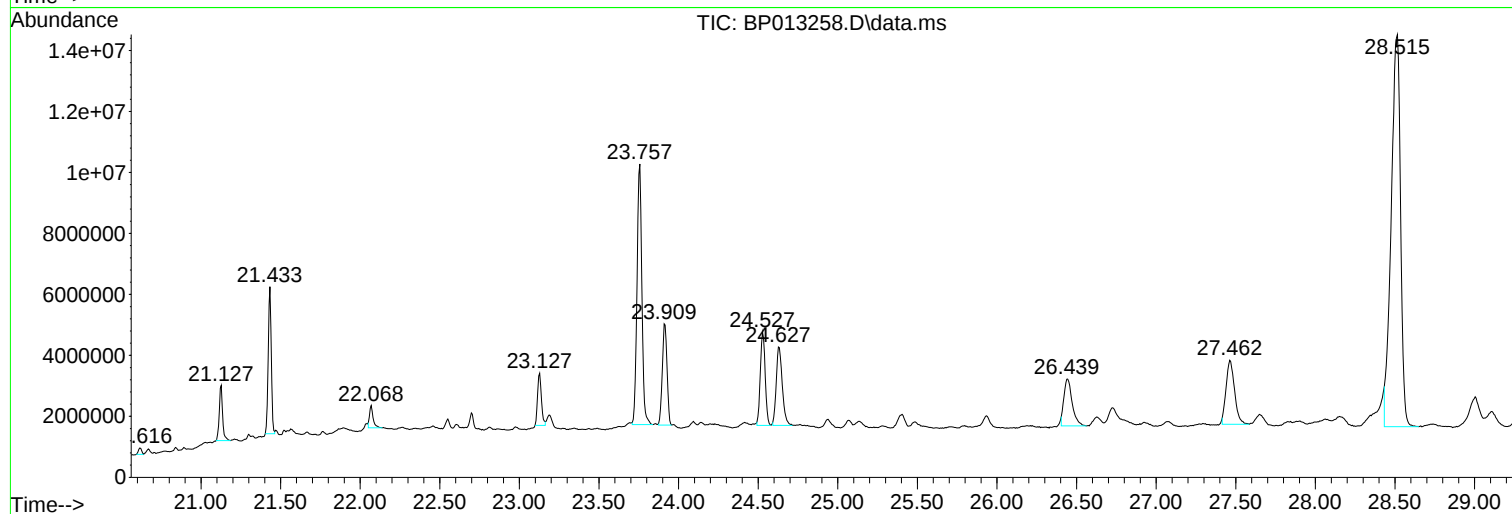
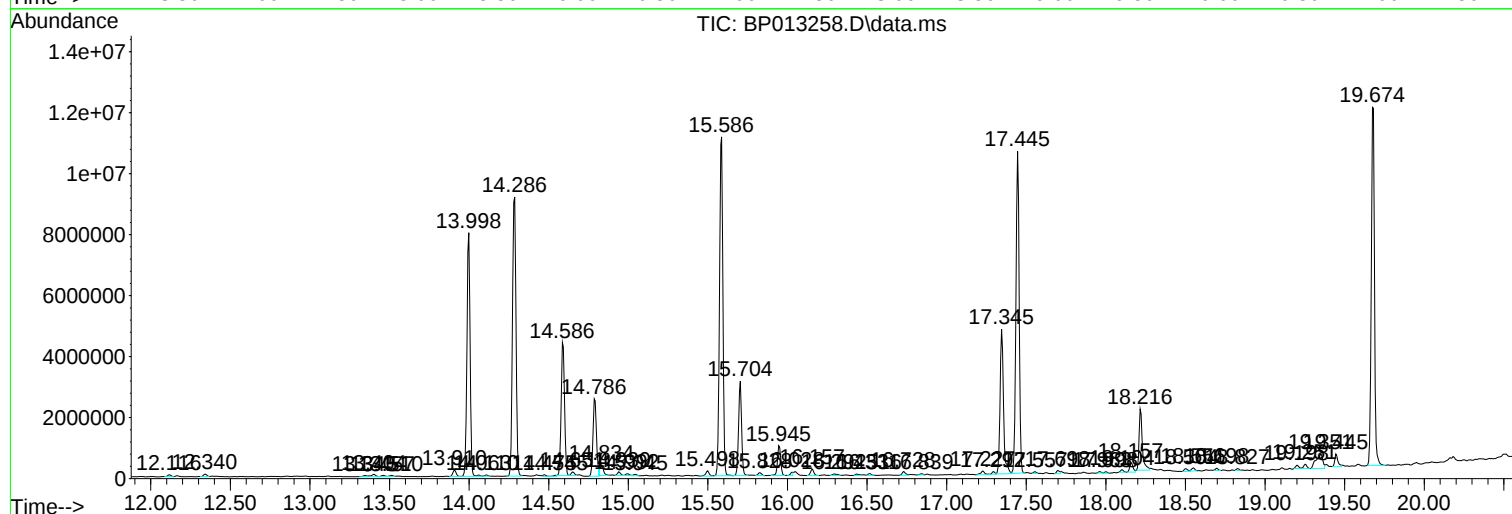
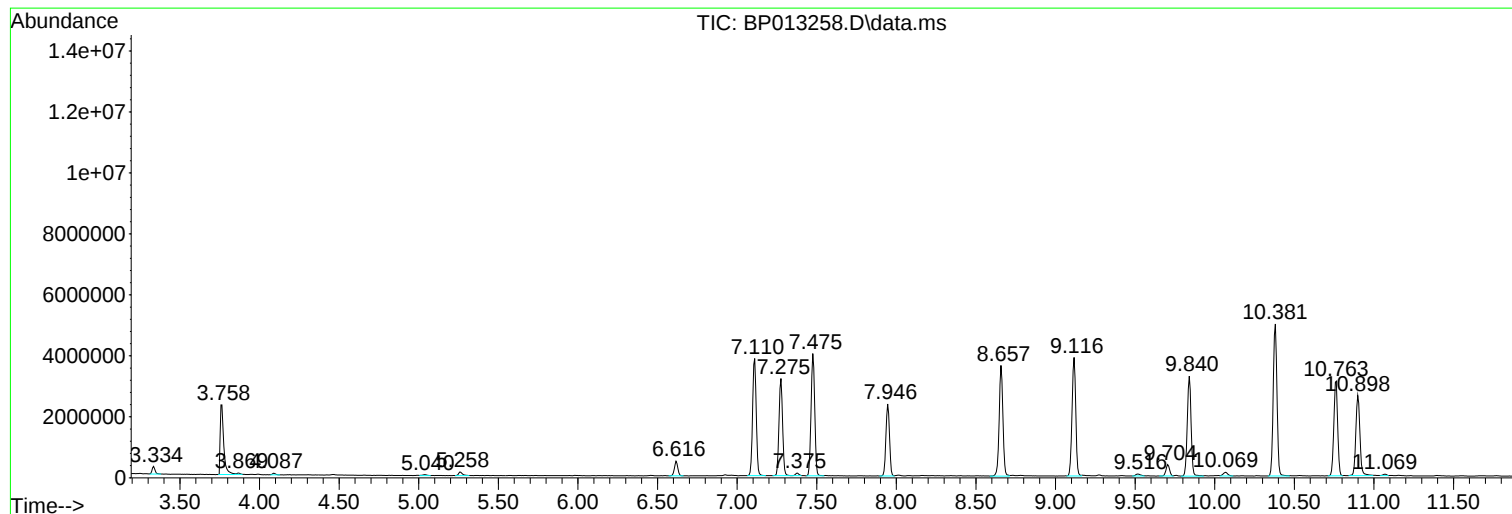
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Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP12222\
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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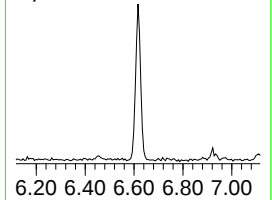
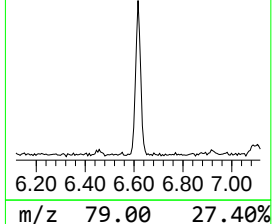
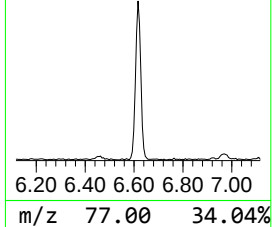
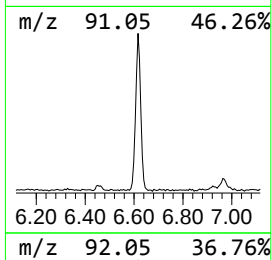
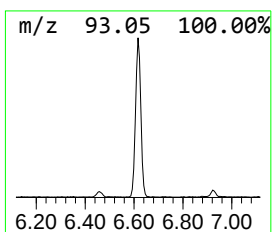
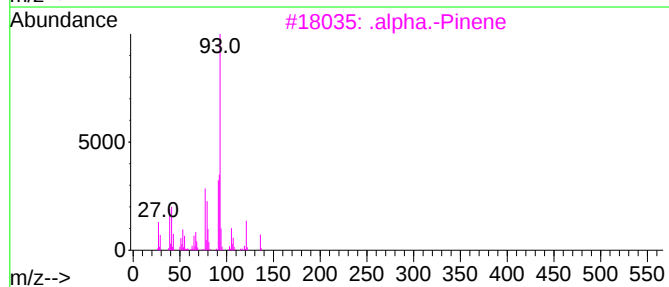
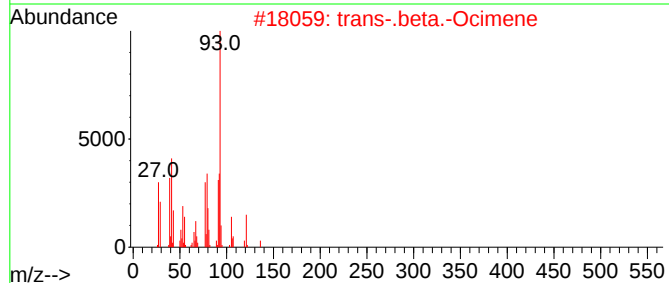
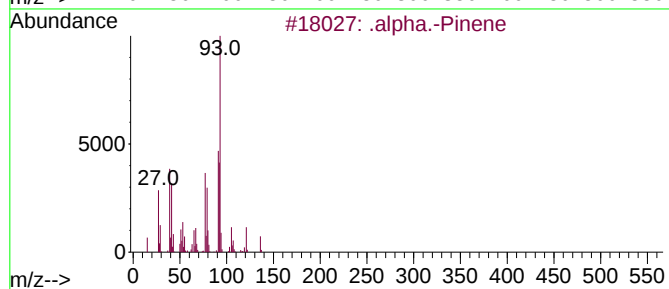
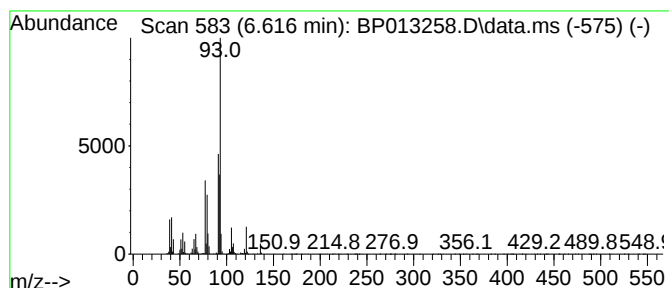
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 .alpha.-Pinene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.616	3.88 ng/ul	709045	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.		.alpha.-Pinene	136	C10H16	000080-56-8	95
2	trans-.beta.-Ocimene			136	C10H16	003779-61-1	94
3	.		.alpha.-Pinene	136	C10H16	000080-56-8	91
4	(+)-3-Carene			136	C10H16	000498-15-7	91
5	Tricyclo[2.2.1.0(2,6)]heptane, 1...			136	C10H16	000508-32-7	91



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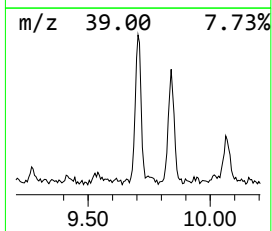
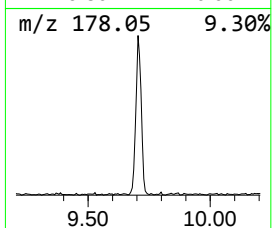
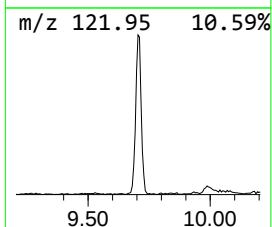
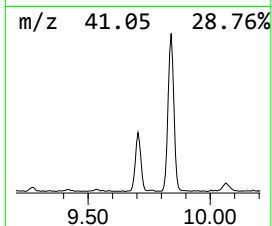
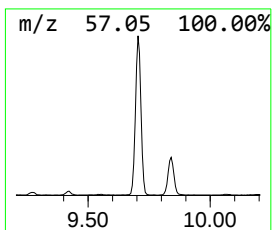
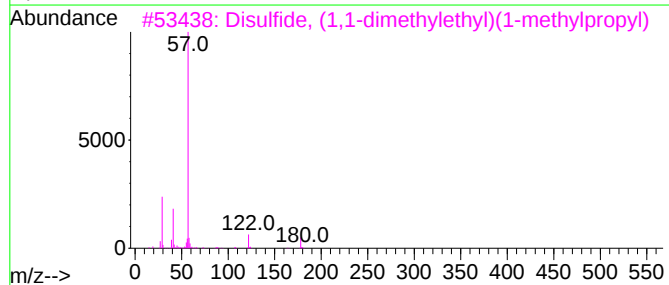
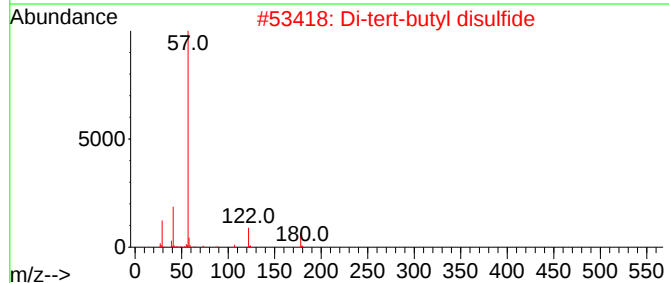
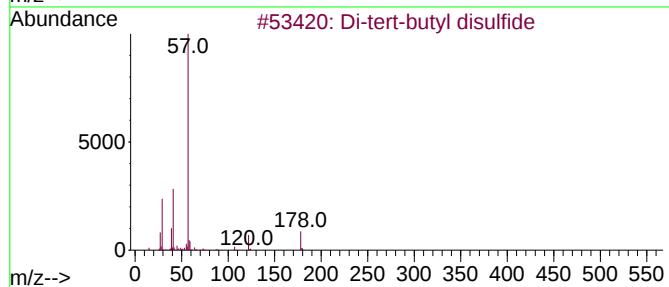
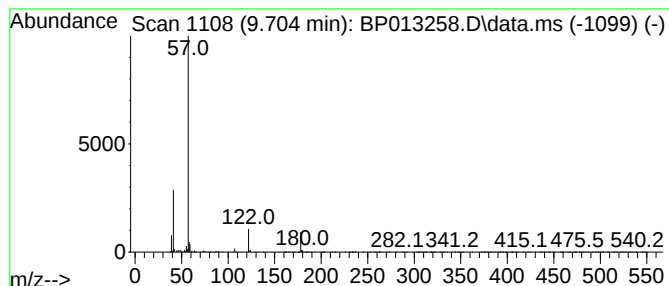
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Di-tert-butyl disulfide Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.704	2.23 ng/ul	580670	Naphthalene-d8	10.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Di-tert-butyl disulfide	178	C8H18S2	000110-06-5	90
2			Di-tert-butyl disulfide	178	C8H18S2	000110-06-5	80
3			Disulfide, (1,1-dimethylethyl)(1...	178	C8H18S2	072437-46-8	80
4			Di-tert-butyl disulfide	178	C8H18S2	000110-06-5	64
5			Disulfide, dibutyl	178	C8H18S2	000629-45-8	58



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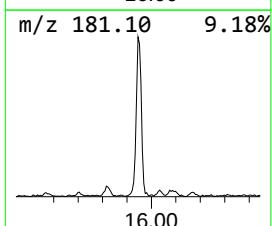
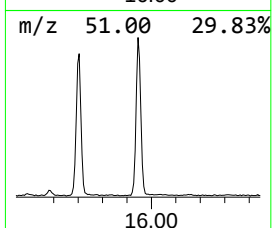
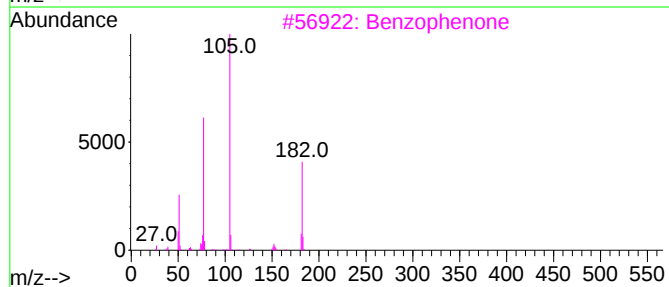
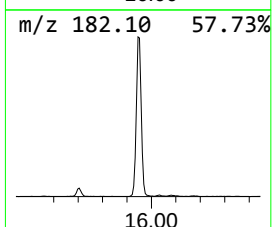
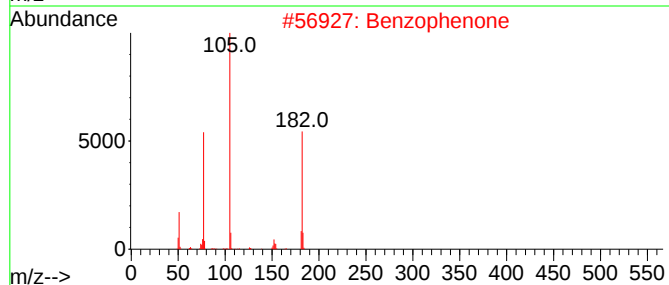
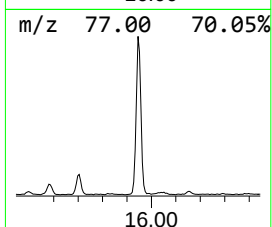
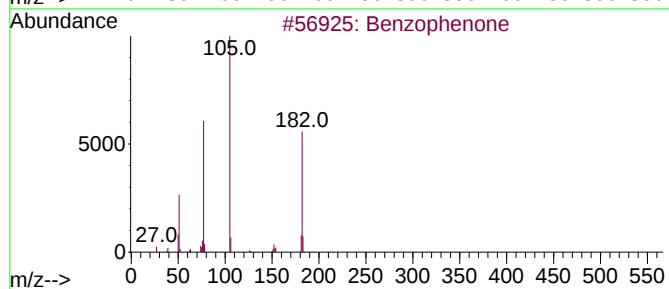
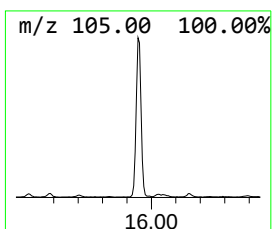
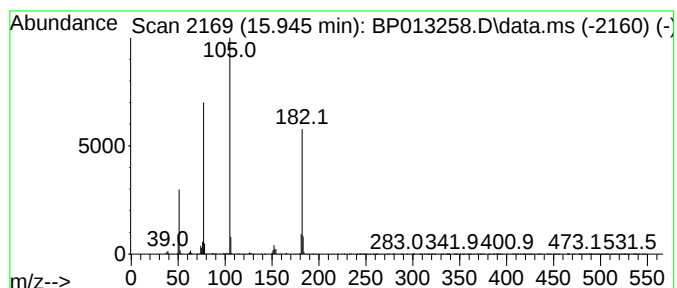
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.945	4.17 ng/ul	1352180	Acenaphthene-d10	14.586

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	91
5			Benzophenone	182	C13H10O	000119-61-9	90



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DBXZ8

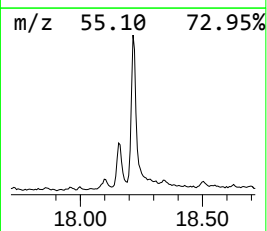
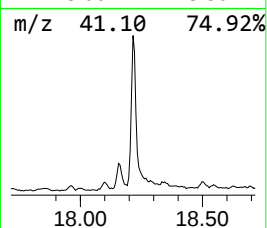
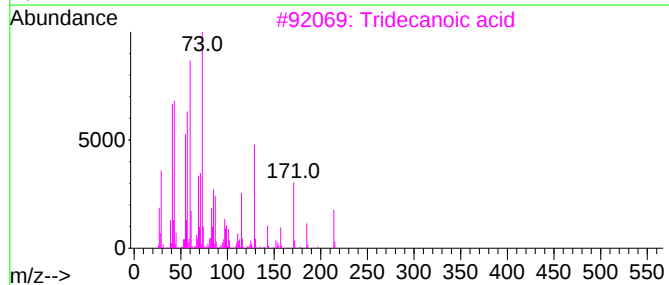
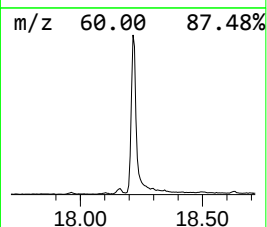
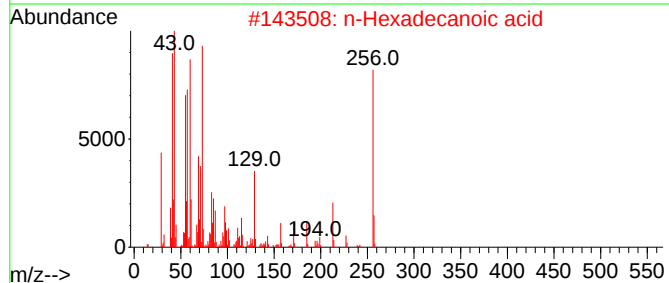
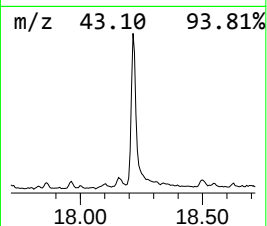
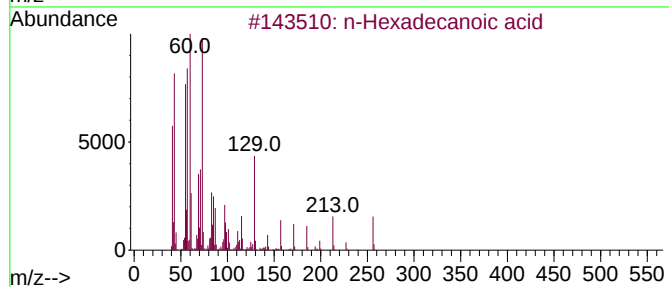
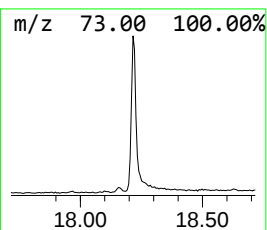
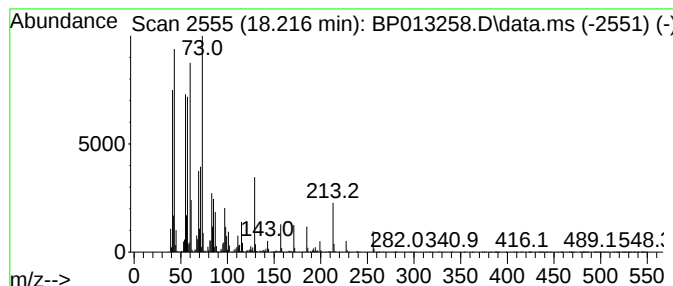
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	7.95 ng/ul	2824410	Phenanthrene-d10	17.345

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		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	95
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013258.D
Acq On : 22 Dec 2022 23:00
Operator : CG/JU
Sample : N6015-19
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ8

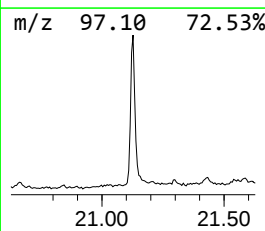
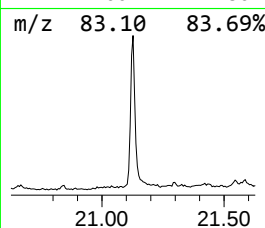
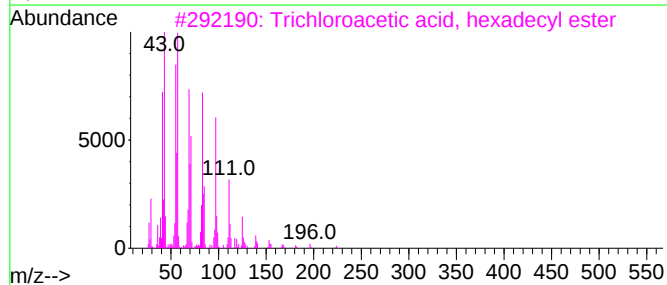
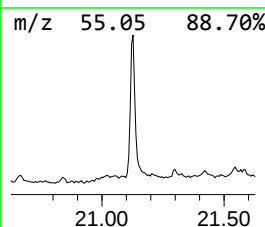
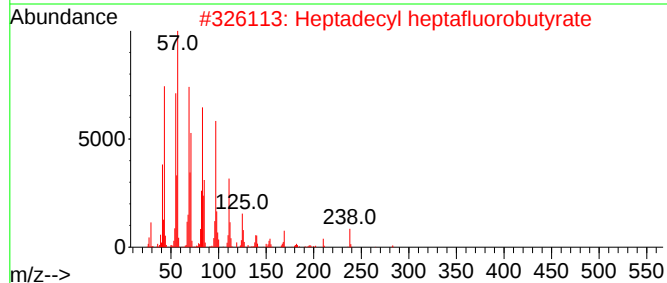
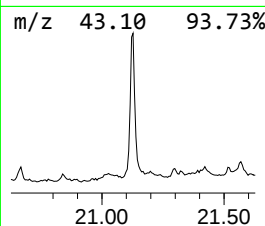
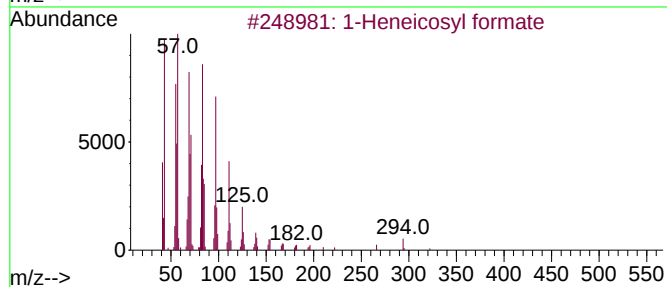
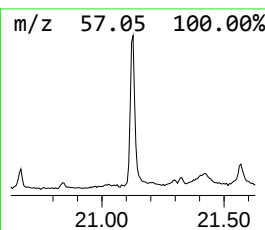
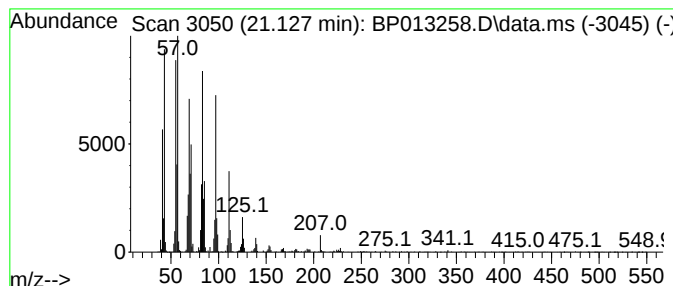
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Heneicosyl formate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.127	8.03 ng/ul	2613050	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	97
2			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
4			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5			9-Tricosene, (Z)-	322	C23H46	027519-02-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013258.D
Acq On : 22 Dec 2022 23:00
Operator : CG/JU
Sample : N6015-19
Misc :
ALS Vial : 25 Sample Multiplier: 1

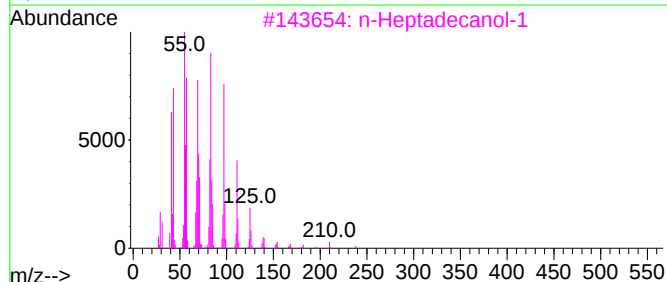
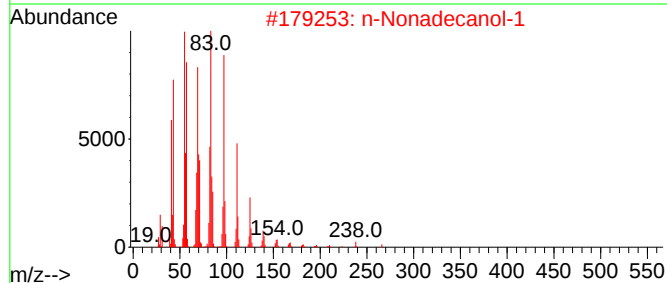
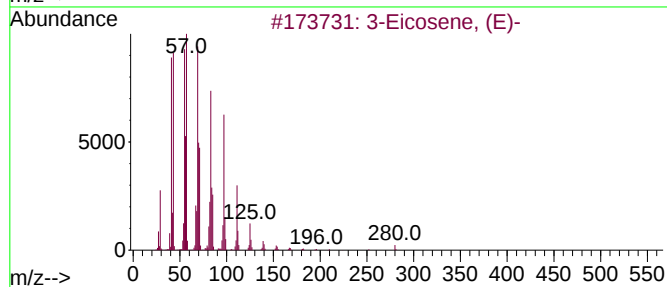
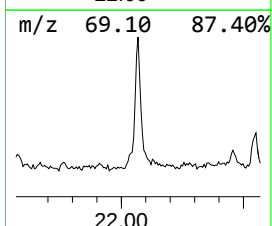
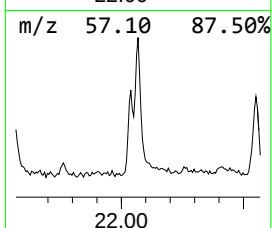
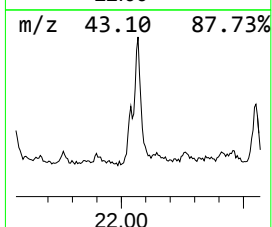
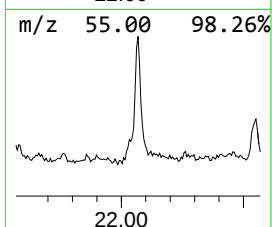
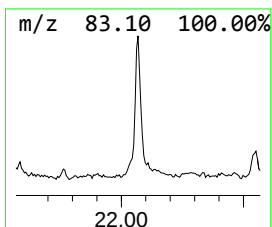
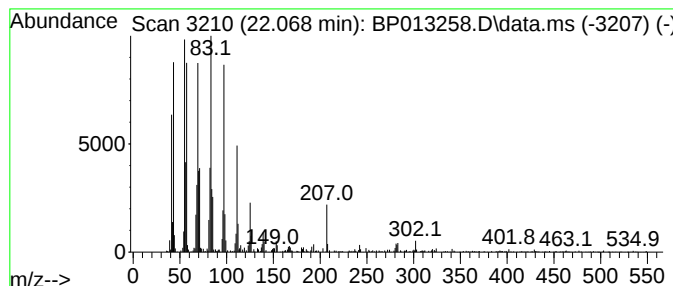
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.068	3.41 ng/ul	1108700	Chrysene-d12	21.433	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	95
2	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
3	n-Heptadecanol-1	256	C17H36O	001454-85-9	94
4	1-Octadecanol	270	C18H38O	000112-92-5	94
5	1-Heneicosanol	312	C21H44O	015594-90-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013258.D
Acq On : 22 Dec 2022 23:00
Operator : CG/JU
Sample : N6015-19
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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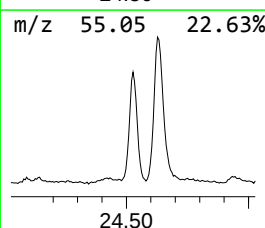
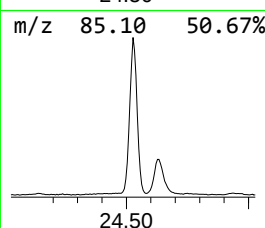
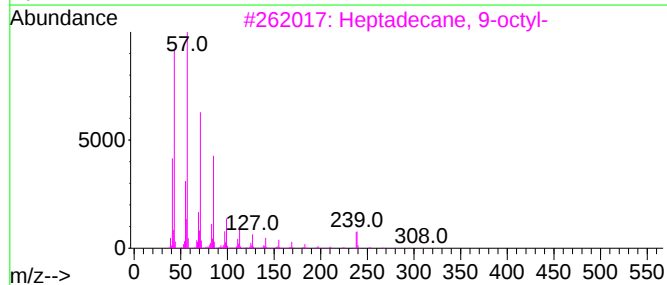
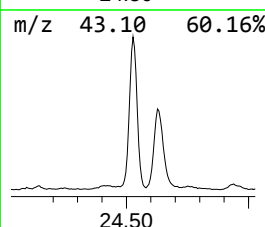
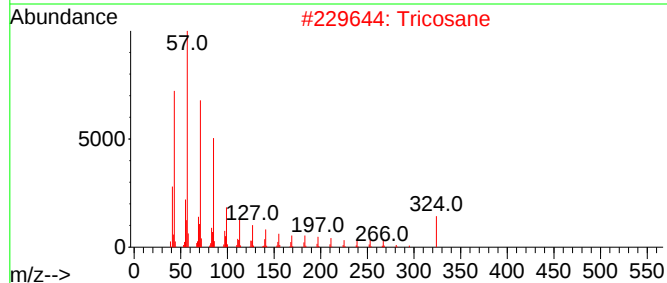
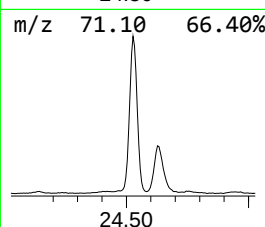
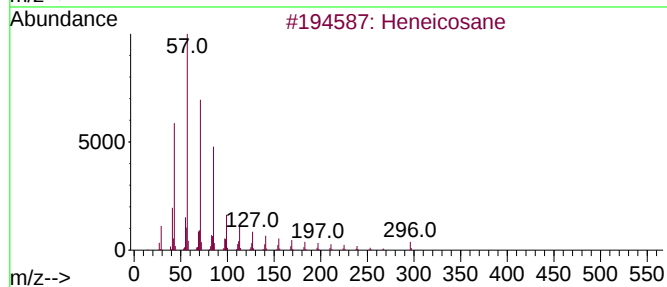
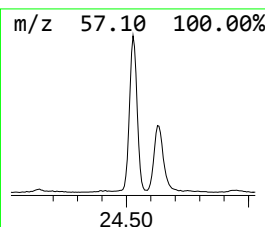
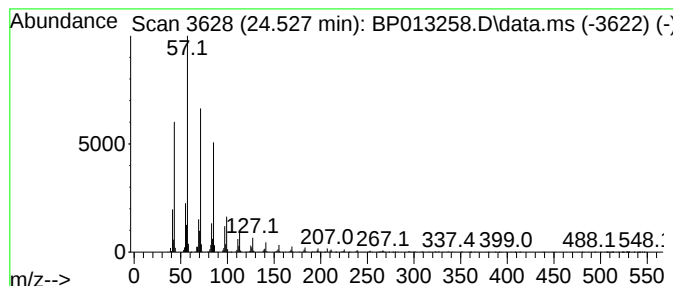
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.527	18.95 ng/ul	6481220	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	97
2			Tricosane	324	C23H48	000638-67-5	95
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
4			Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
5			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013258.D
Acq On : 22 Dec 2022 23:00
Operator : CG/JU
Sample : N6015-19
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ8

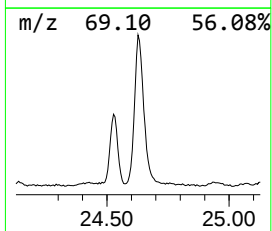
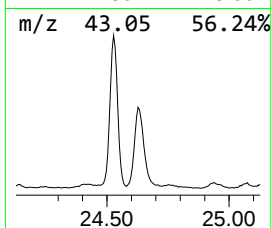
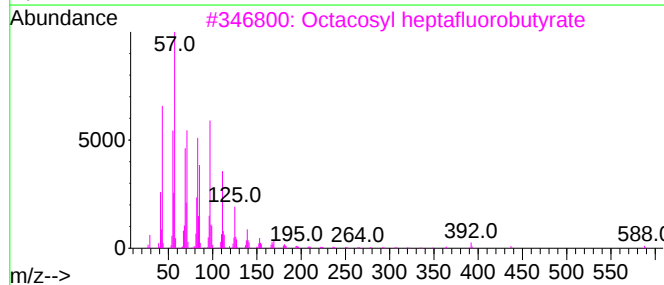
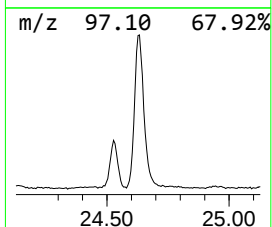
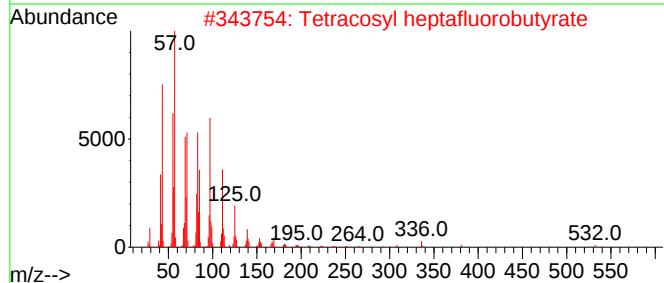
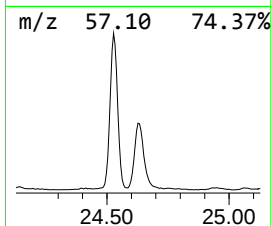
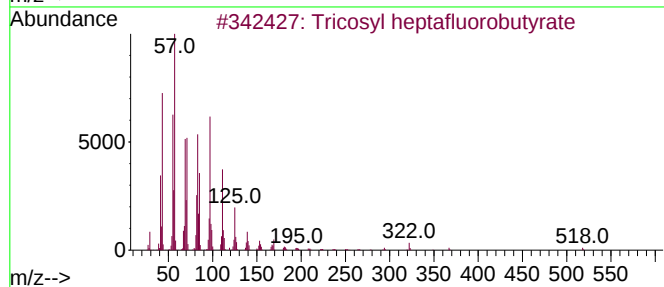
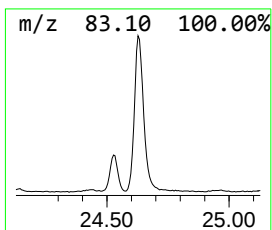
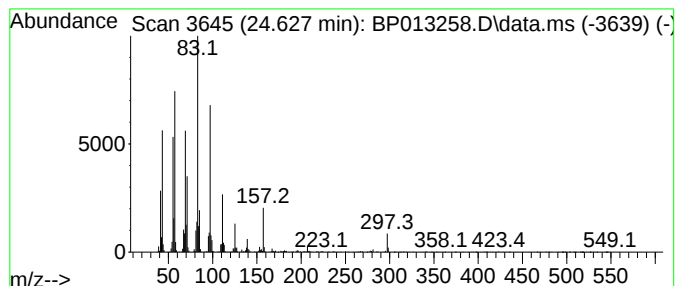
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.627	19.79 ng/ul	6769380	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricosyl heptafluorobutyrate	536	C27H47F7O2	1000351-83-4	49
2			Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	49
3			Octacosyl heptafluorobutyrate	606	C32H57F7O2	1010351-83-6	49
4			Hexacosyl pentafluoropropionate	528	C29H53F5O2	1000351-81-2	46
5			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013258.D
Acq On : 22 Dec 2022 23:00
Operator : CG/JU
Sample : N6015-19
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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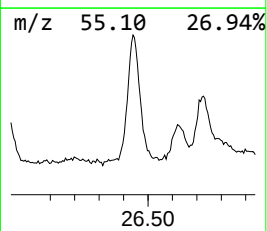
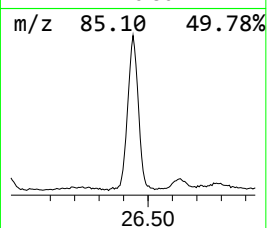
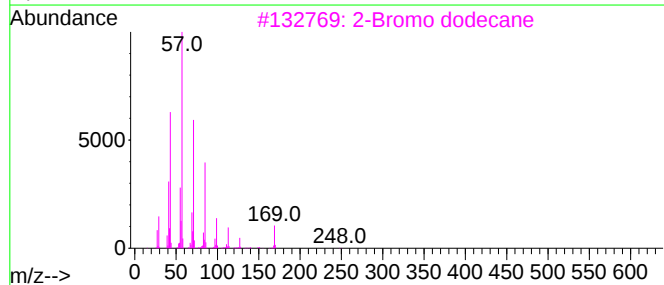
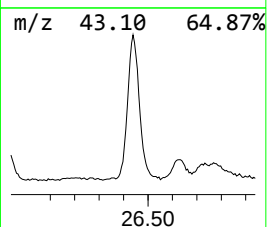
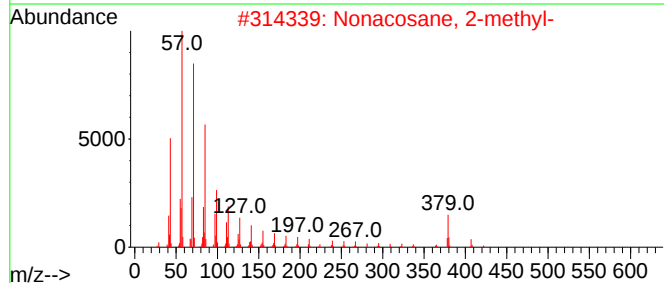
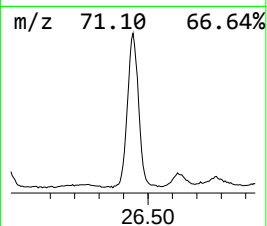
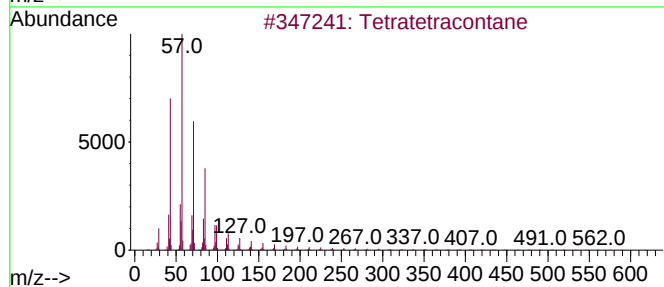
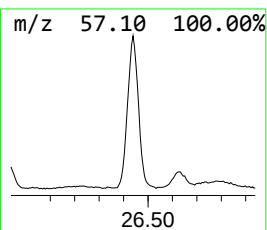
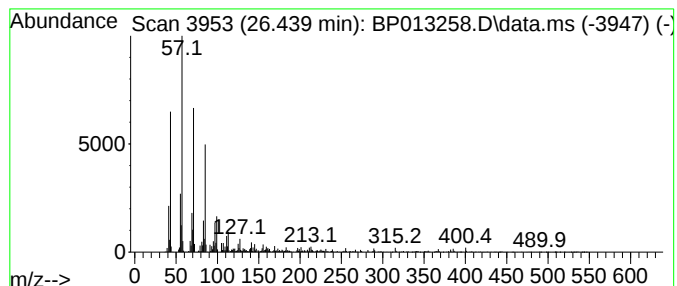
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.439	16.19 ng/ul	5535990	Perylene-d12	23.910

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetratetracontane	619	C44H90	007098-22-8	94
2		Nonacosane, 2-methyl-	422	C30H62	001560-75-4	94
3		2-Bromo dodecane	248	C12H25Br	013187-99-0	93
4		Eicosane	282	C20H42	000112-95-8	91
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013258.D
Acq On : 22 Dec 2022 23:00
Operator : CG/JU
Sample : N6015-19
Misc :
ALS Vial : 25 Sample Multiplier: 1

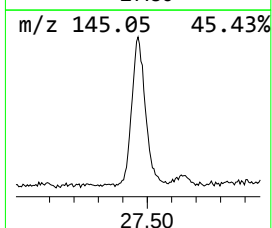
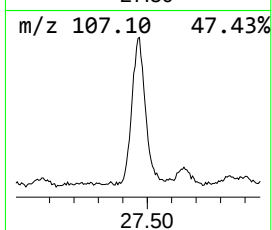
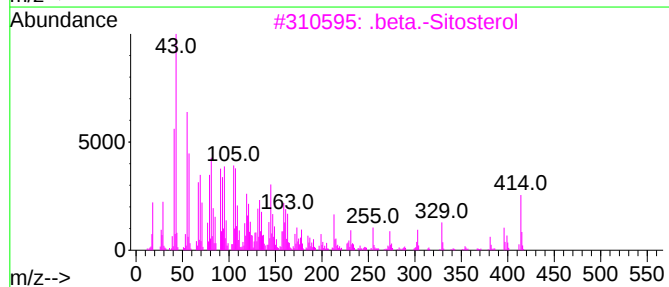
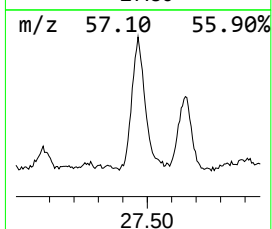
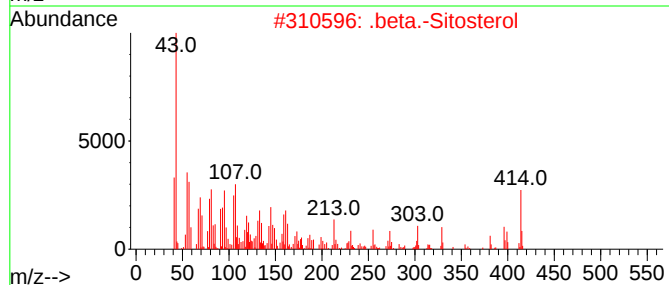
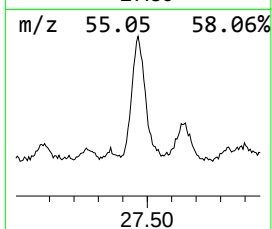
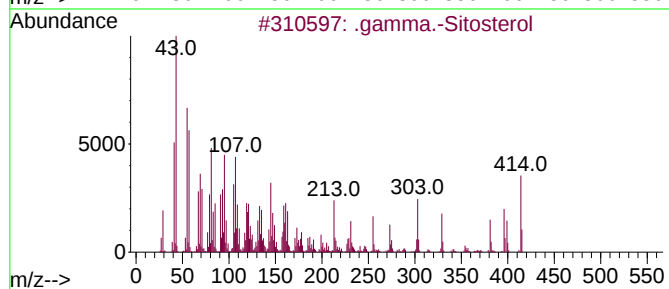
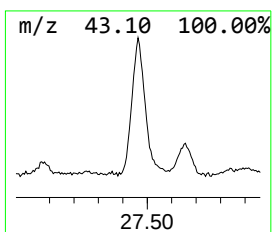
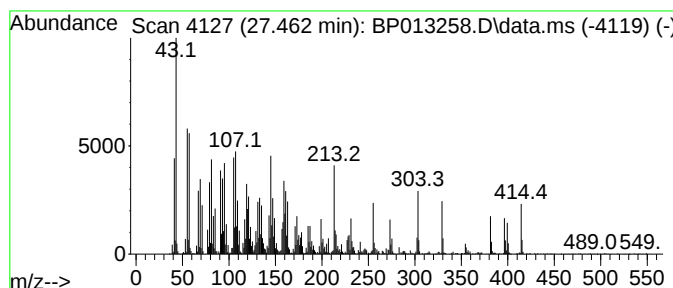
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 .gamma.-Sitosterol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
27.462	22.58 ng/ul	7722900	Perylene-d12		23.910	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol		414	C29H50O	000083-47-6	99
2	.beta.-Sitosterol		414	C29H50O	000083-46-5	94
3	.beta.-Sitosterol		414	C29H50O	000083-46-5	93
4	.gamma.-Sitosterol		414	C29H50O	000083-47-6	53
5	Campesterol		400	C28H48O	000474-62-4	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013258.D
Acq On : 22 Dec 2022 23:00
Operator : CG/JU
Sample : N6015-19
Misc :
ALS Vial : 25 Sample Multiplier: 1

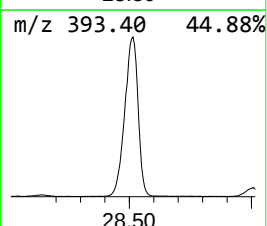
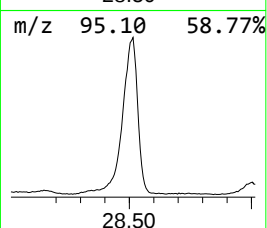
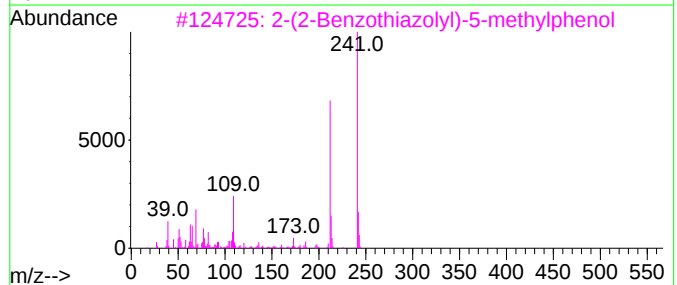
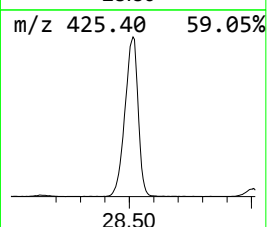
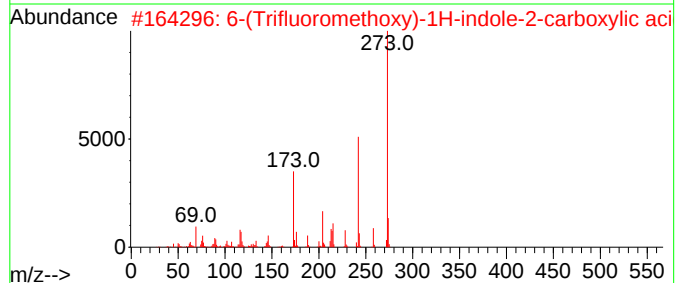
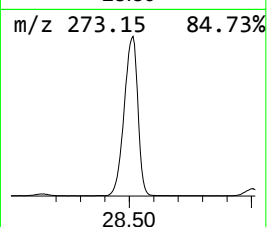
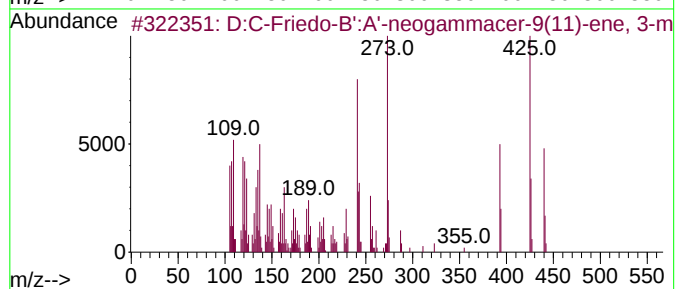
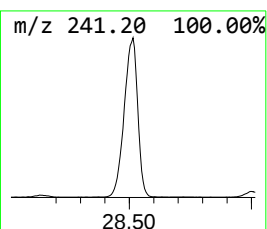
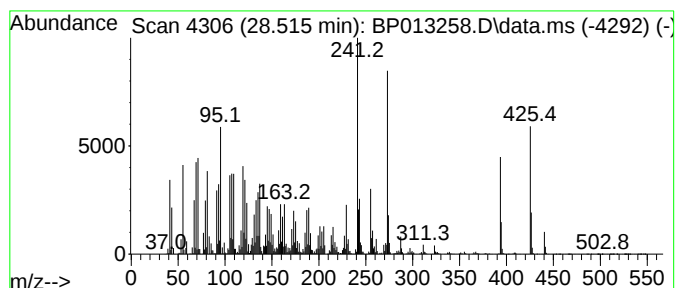
Instrument :
BNA_P
ClientSampleId :
DBXZ8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 D:C-Friedo-B':A'-neogammace... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.515	153.69 ng/ul	52563800	Perylene-d12	23.910
Hit#	of 5	Tentative ID	MW MolForm	CAS# Qual
1		D:C-Friedo-B':A'-neogammacer-9(1...	440 C31H52O	004555-56-0 50
2		6-(Trifluoromethoxy)-1H-indole-2...	273 C12H10F3NO3	1000499-75-7 25
3		2-(2-Benzothiazolyl)-5-methylphenol	241 C14H11NOS	056048-54-5 15
4		Phenol, 2,6-dimethyl-, phosphite...	394 C24H27O3P	052830-49-6 15
5		Pyrazolo[1,5-a]pyrimidine, 6,7-d...	273 C18H15N3	186956-30-9 15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013258.D
Acq On : 22 Dec 2022 23:00
Operator : CG/JU
Sample : N6015-19
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA 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
.alpha.-Pinene	6.616	3.9	ng/ul	709045	1	7.946	3653490	20.0
Di-tert-butyl d...	9.704	2.2	ng/ul	580670	2	10.763	5197490	20.0
Benzophenone	15.945	4.2	ng/ul	1352180	3	14.586	6482300	20.0
n-Hexadecanoic ...	18.216	8.0	ng/ul	2824410	4	17.345	7104490	20.0
1-Heneicosyl fo...	21.127	8.0	ng/ul	2613050	5	21.433	6511920	20.0
(DEL) Alkane: S...	22.068	3.4	ng/ul	1108700	5	21.433	6511920	20.0
(DEL) Alkane: S...	24.527	18.9	ng/ul	6481220	6	23.910	6840480	20.0
unknown-01	24.627	19.8	ng/ul	6769380	6	23.910	6840480	20.0
(DEL) Alkane: S...	26.439	16.2	ng/ul	5535990	6	23.910	6840480	20.0
.gamma.-Sitosterol	27.462	22.6	ng/ul	7722900	6	23.910	6840480	20.0
D:C-Friedo-B':A...	28.515	153.7	ng/ul	52563800	6	23.910	6840480	20.0