

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062723\
 Data File : BP015752.D
 Acq On : 27 Jun 2023 18:03
 Operator : MA/JU
 Sample : 03085-17
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFQ3

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-BP062623.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015752.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.058	142	148	155	rVB	18748	32158	0.66%	0.117%
2	4.158	161	165	168	rBV2	10484	14496	0.30%	0.053%
3	5.111	324	327	332	rVB7	4791	7202	0.15%	0.026%
4	7.240	682	689	706	rBV	240563	603472	12.39%	2.203%
5	7.387	709	714	731	rVB	227596	426911	8.77%	1.558%
6	7.593	741	749	764	rBV	337700	643730	13.22%	2.350%
7	7.940	805	808	813	rVB7	3677	5851	0.12%	0.021%
8	8.063	821	829	849	rBV	429803	802226	16.47%	2.929%
9	8.269	859	864	867	rBV7	3544	5563	0.11%	0.020%
10	8.793	945	953	969	rVV	298900	700653	14.39%	2.558%
11	8.916	969	974	986	rVV2	36454	88357	1.81%	0.323%
12	9.252	1024	1031	1047	rBV	261997	553030	11.36%	2.019%
13	9.981	1148	1155	1173	rBV	213788	490720	10.08%	1.791%
14	10.546	1243	1251	1269	rBV	187488	634124	13.02%	2.315%
15	10.893	1302	1310	1321	rBV	523917	968193	19.88%	3.535%
16	11.063	1331	1339	1362	rBV	176675	609823	12.52%	2.226%
17	13.592	1765	1769	1775	rVB4	13231	22176	0.46%	0.081%
18	14.122	1853	1859	1872	rBV	510934	922032	18.93%	3.366%
19	14.410	1901	1908	1923	rBV	708408	1139152	23.39%	4.159%
20	14.716	1953	1960	1971	rBV	700271	1131806	23.24%	4.132%
21	14.998	2002	2008	2024	rBV3	63812	307728	6.32%	1.123%
22	15.710	2123	2129	2143	rBV	715734	1367571	28.08%	4.993%
23	15.875	2152	2157	2168	rBV2	114072	277220	5.69%	1.012%
24	17.475	2423	2429	2440	rBV	824791	1330272	27.31%	4.856%
25	17.575	2440	2446	2463	rVB2	779684	1408332	28.92%	5.141%
26	18.786	2646	2652	2655	rBV5	68987	166853	3.43%	0.609%
27	19.804	2819	2825	2835	rBV	1168713	2289378	47.01%	8.358%
28	21.563	3120	3124	3130	rBV	1024874	1744508	35.82%	6.369%
29	21.663	3134	3141	3170	rVB6	923083	4870228	100.00%	17.779%
30	23.939	3522	3528	3540	rVB2	922131	2156196	44.27%	7.871%
31	24.110	3551	3557	3565	rBV	714073	1672545	34.34%	6.106%

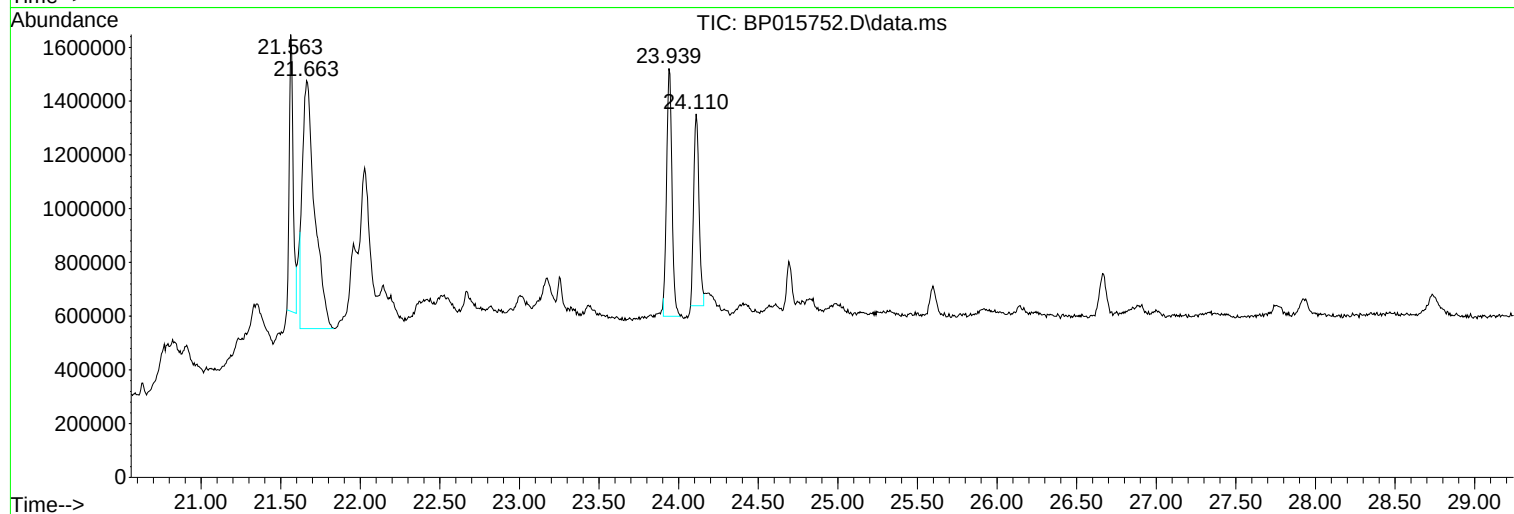
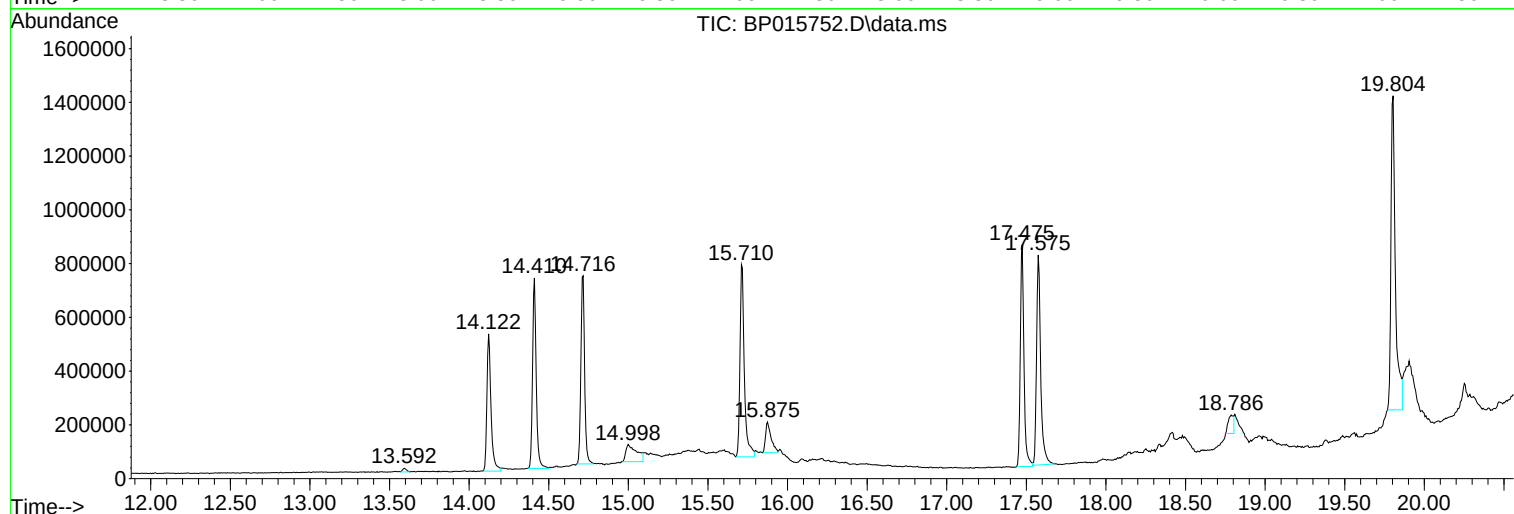
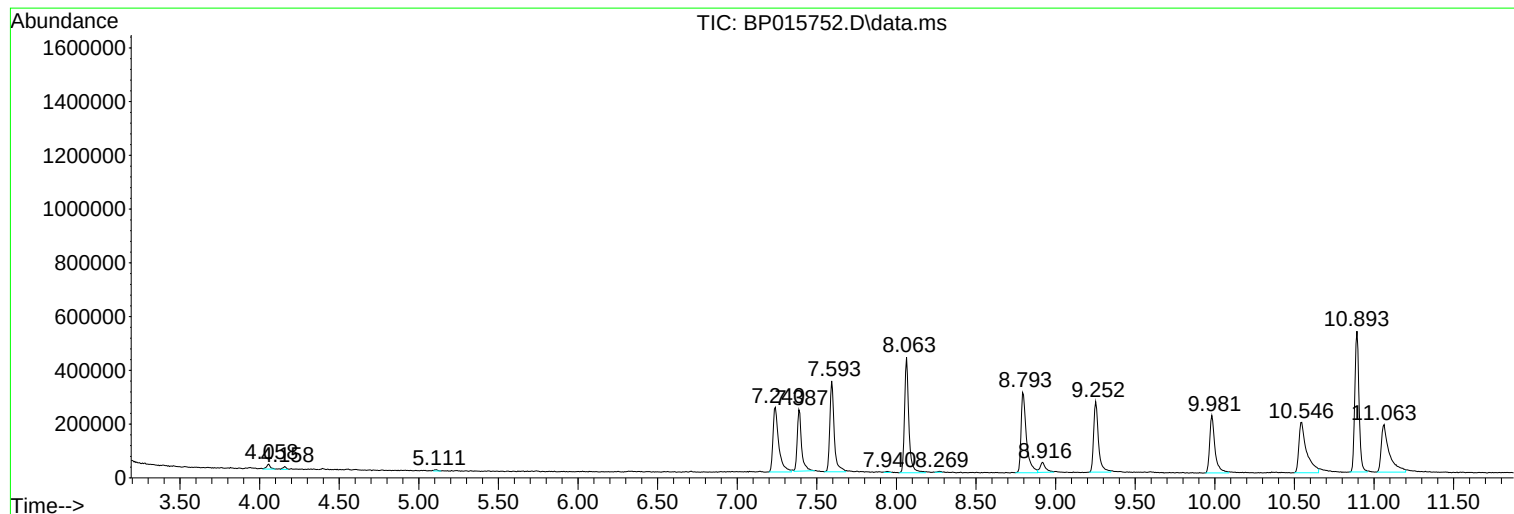
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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



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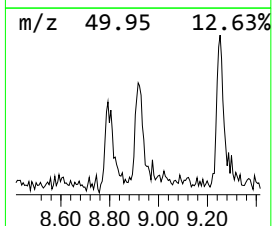
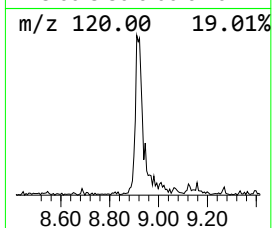
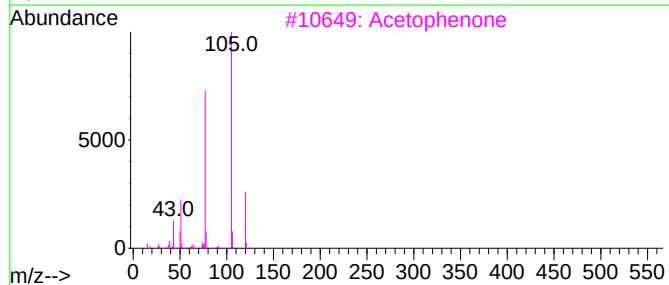
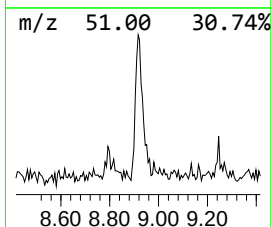
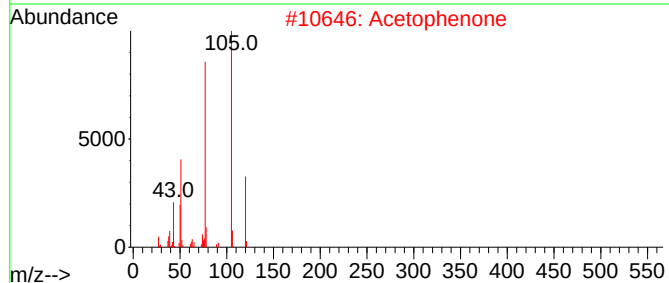
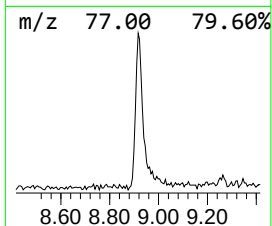
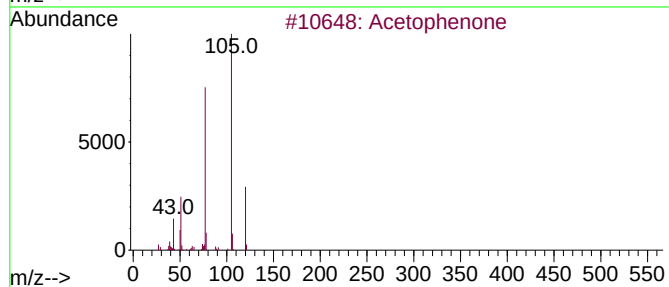
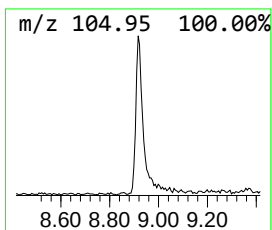
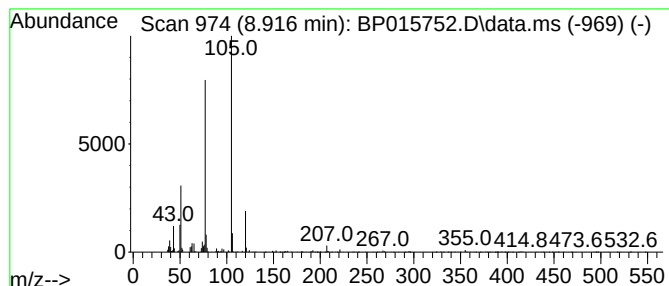
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Aceto phenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.916	2.20 ng/ul	88357	1,4-Dichlorobenzene-d4	8.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone			120	C8H8O	000098-86-2	91
2	Acetophenone			120	C8H8O	000098-86-2	91
3	Acetophenone			120	C8H8O	000098-86-2	91
4	Acetophenone			120	C8H8O	000098-86-2	90
5	Acetophenone			120	C8H8O	000098-86-2	90



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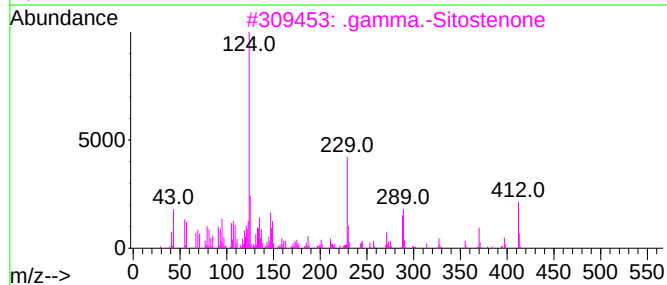
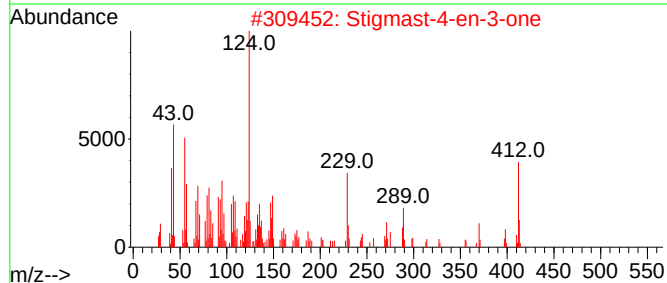
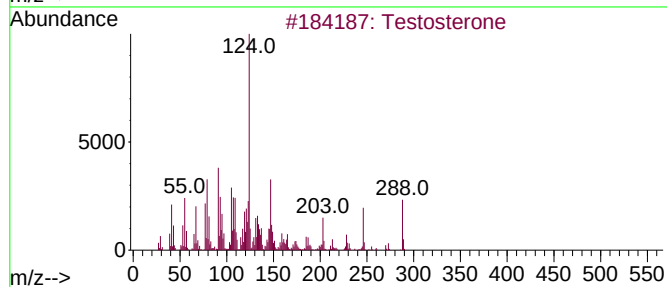
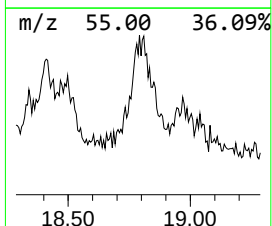
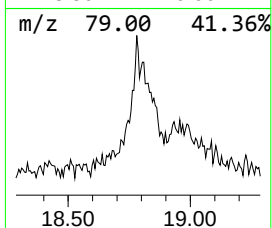
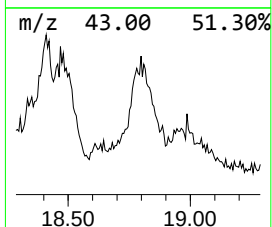
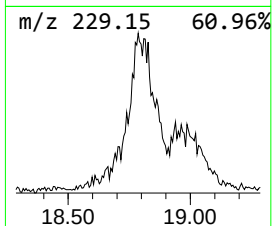
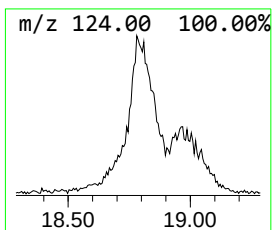
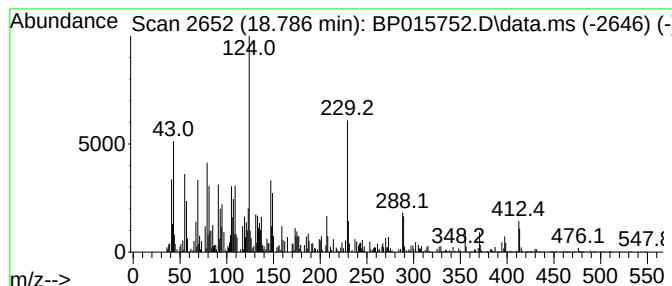
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Testosterone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.786	2.51 ng/ul	166853	Phenanthrene-d10	17.475

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Testosterone	288	C19H28O2	000058-22-0	87
2			Stigmast-4-en-3-one	412	C29H48O	001058-61-3	87
3			.gamma.-Sitostenone	412	C29H48O	084924-96-9	87
4			Testosterone	288	C19H28O2	000058-22-0	50
5			Testosterone cypionate	412	C27H40O3	000058-20-8	49



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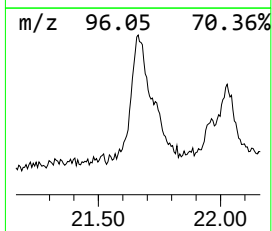
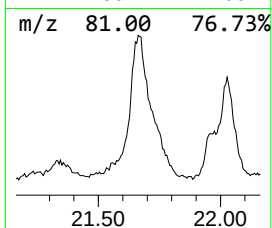
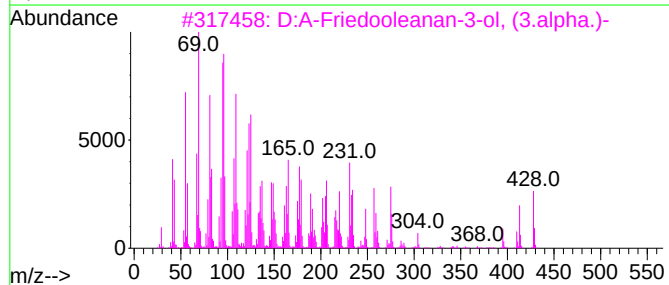
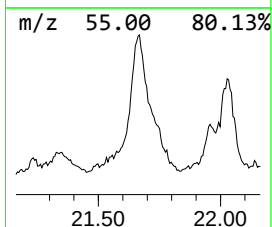
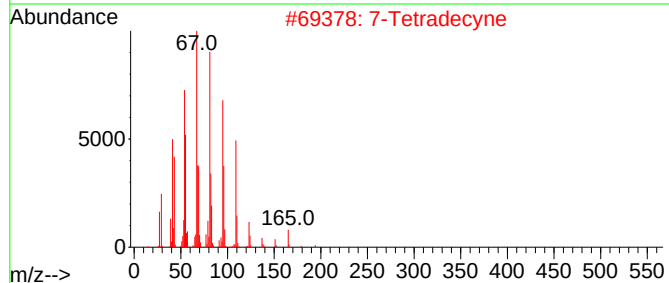
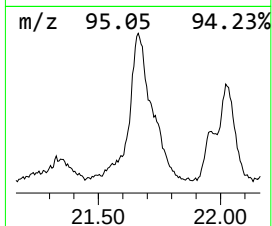
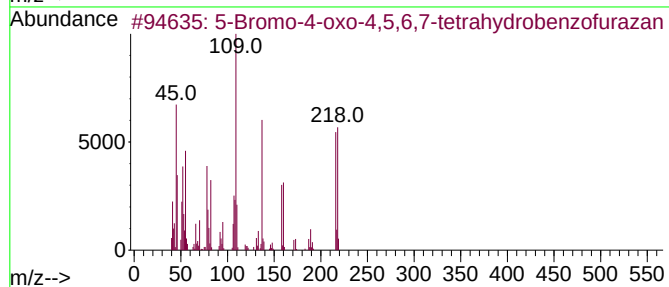
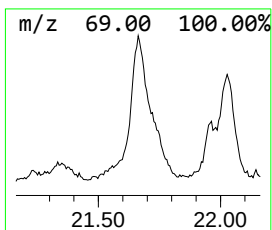
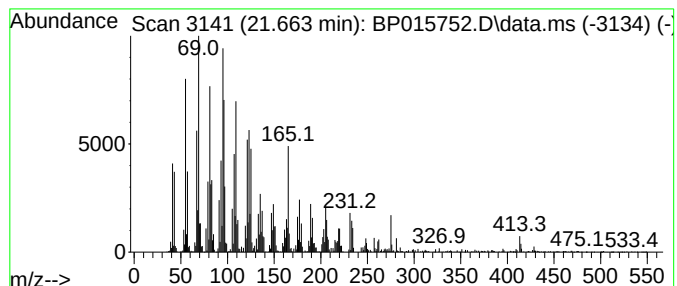
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.663	55.84 ng/ul	4870230	Chrysene-d12	21.563

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	59
2		7-Tetradecyne	194	C14H26	035216-11-6	55
3		D:A-Friedooleanan-3-ol, (3.alpha.)-	428	C30H52O	005085-72-3	53
4		Naphthalene, decahydro-1,1-dimet...	166	C12H22	035431-04-0	46
5		13-Tetradecene-11-yn-1-ol	208	C14H24O	1000131-00-4	45



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Aceto phenone	8.916	2.2	ng/u1	88357	1	8.063	802226	20.0
Testosterone	18.786	2.5	ng/u1	166853	4	17.475	1330270	20.0
5-Bromo-4-oxo-4...	21.663	55.8	ng/u1	4870230	5	21.563	1744510	20.0