

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081318\
 Data File : BF108116.D
 Acq On : 13 Aug 2018 15:16
 Operator : JU/SJ
 Sample : J4272-19 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-03-080918-J

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % 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_F\METHODS\8270-BF080818.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.254	491	496	502	rBV	606728	647982	19.09%	1.818%
2	5.643	558	562	565	rBV	3601885	3194269	94.08%	8.964%
3	6.637	726	731	741	rBV	3815102	3395109	100.00%	9.527%
4	6.790	753	757	760	rBV	3879201	3305118	97.35%	9.275%
5	7.025	793	797	800	rBV	1828231	1590691	46.85%	4.464%
6	7.178	819	823	826	rBV	3092412	2467212	72.67%	6.923%
7	7.578	887	891	895	rBV	2202933	2074902	61.11%	5.823%
8	8.307	1011	1015	1018	rBV	2288056	1908491	56.21%	5.356%
9	9.378	1193	1197	1201	rBV	3594942	3088936	90.98%	8.668%
10	9.772	1261	1264	1267	rBV	254923	195645	5.76%	0.549%
11	10.072	1311	1315	1319	rBV	2084064	1778044	52.37%	4.989%
12	10.866	1446	1450	1463	rBV	1817814	1665941	49.07%	4.675%
13	10.960	1463	1466	1476	rVB4	41927	80526	2.37%	0.226%
14	11.572	1566	1570	1572	rBV	1950663	1645442	48.47%	4.617%
15	11.595	1572	1574	1577	rVV	262730	211123	6.22%	0.592%
16	11.648	1580	1583	1586	rVB	87077	78361	2.31%	0.220%
17	11.907	1624	1627	1631	rVB5	27224	37184	1.10%	0.104%
18	12.072	1652	1655	1658	rBV	90634	91391	2.69%	0.256%
19	12.101	1658	1660	1664	rVB2	67062	69736	2.05%	0.196%
20	12.189	1671	1675	1677	rBV2	86684	91543	2.70%	0.257%
21	12.630	1746	1750	1752	rBV3	104989	123613	3.64%	0.347%
22	12.654	1752	1754	1758	rVB4	66902	77466	2.28%	0.217%
23	12.795	1774	1778	1781	rBV	596620	562998	16.58%	1.580%
24	13.024	1814	1817	1823	rVB	575394	520817	15.34%	1.461%
25	13.160	1836	1840	1843	rBV	3042703	2575565	75.86%	7.227%
26	13.348	1870	1872	1875	rVB	75562	68275	2.01%	0.192%
27	14.230	2017	2022	2025	rBV2	2078512	2185024	64.36%	6.132%
28	14.254	2025	2026	2032	rVB	305315	245481	7.23%	0.689%
29	15.089	2165	2168	2172	rVB	325766	370305	10.91%	1.039%
30	15.807	2286	2290	2294	rBV	974746	1288604	37.95%	3.616%

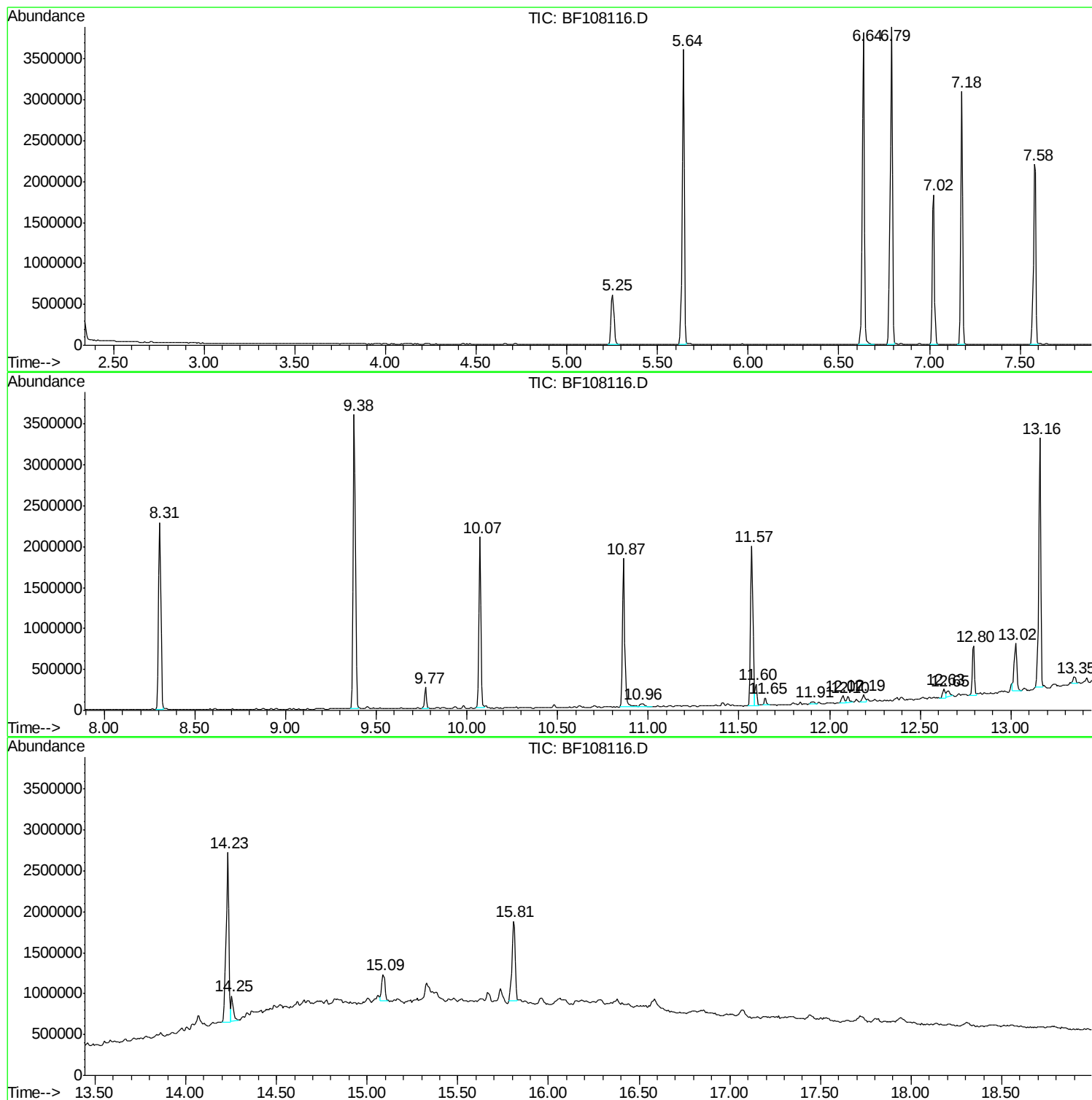
Sum of corrected areas: 35635794

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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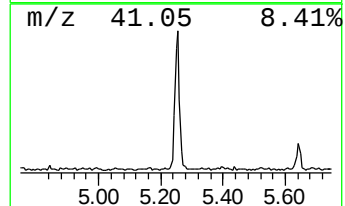
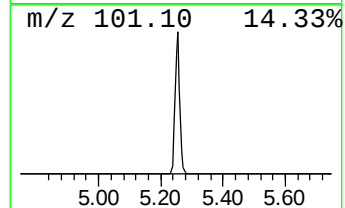
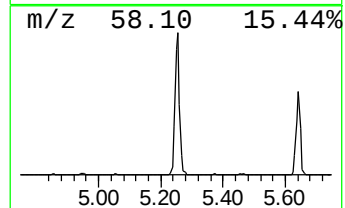
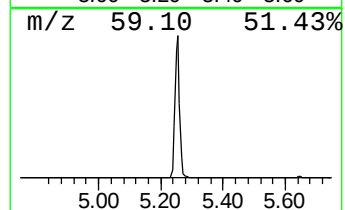
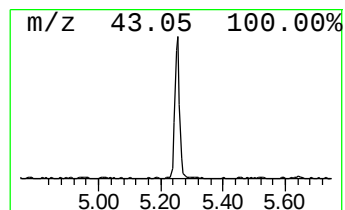
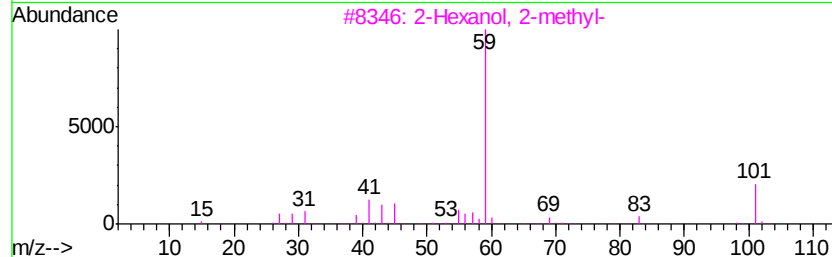
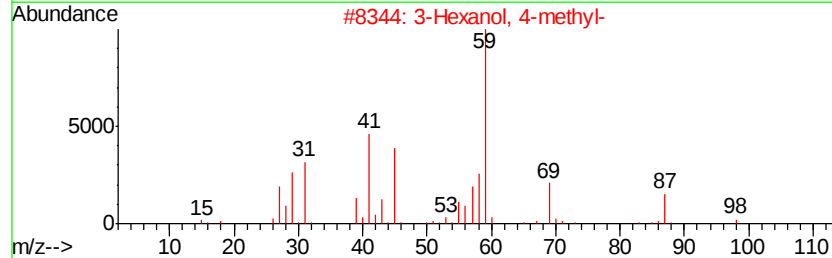
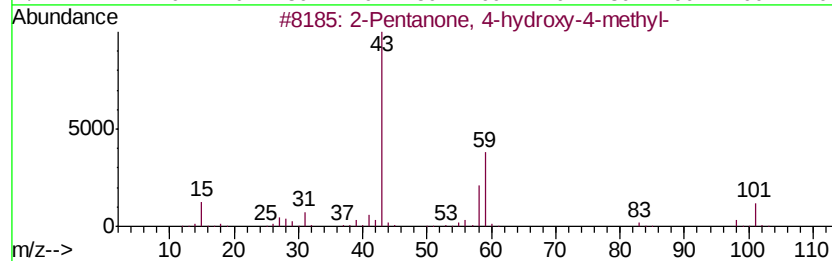
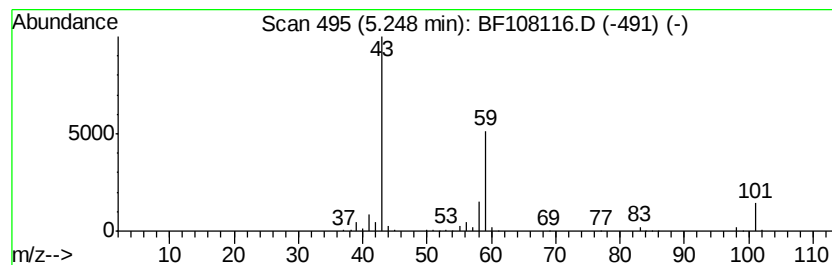
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.25	8.15 ng	647982	1,4-Dichlorobenzene-d4	7.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



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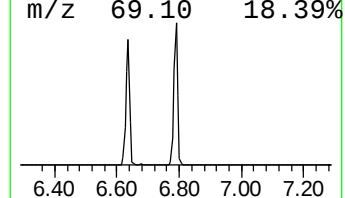
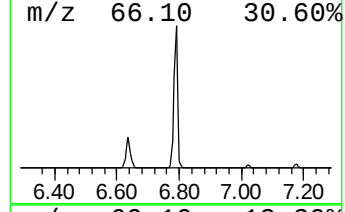
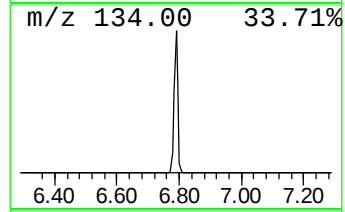
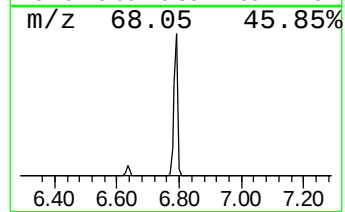
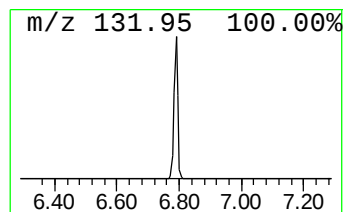
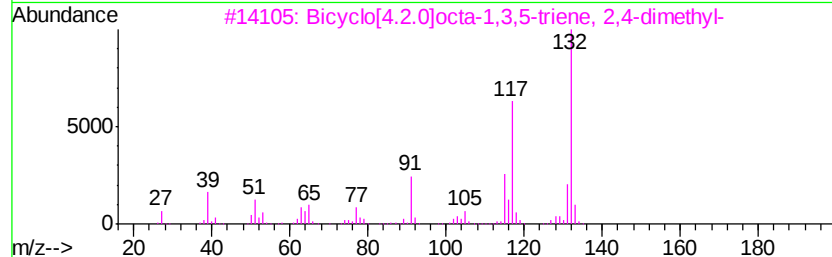
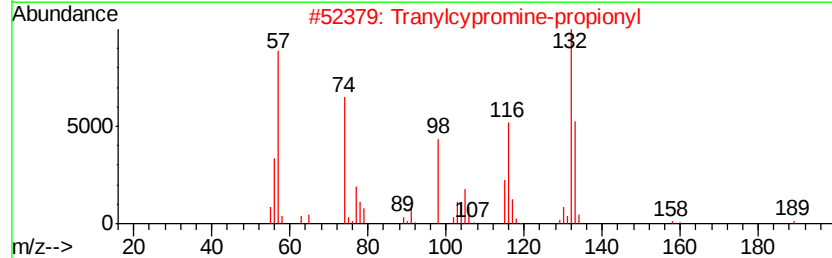
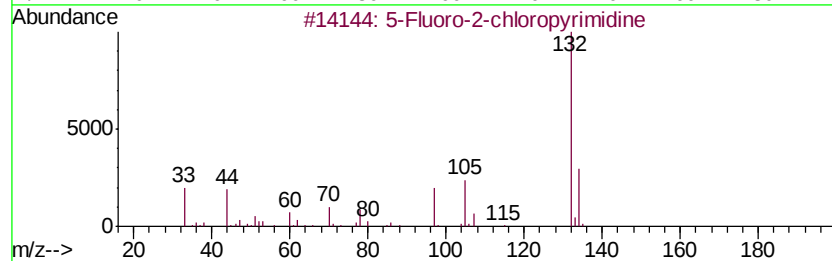
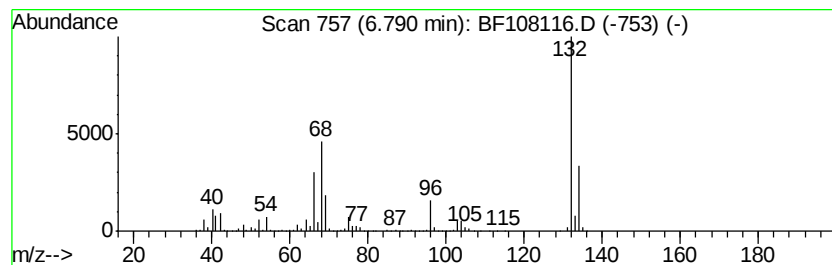
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Peak Number 2 unknown6.79 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.79	41.56 ng	3305120	1,4-Dichlorobenzene-d4	7.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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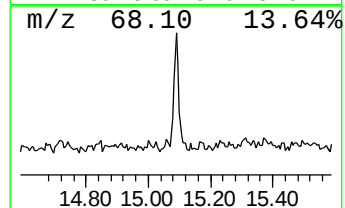
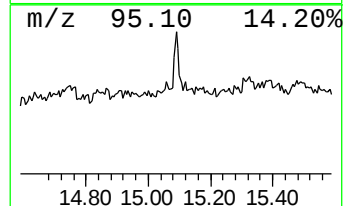
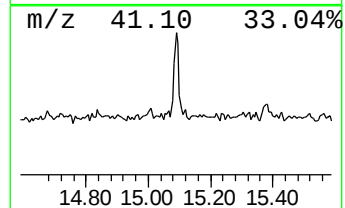
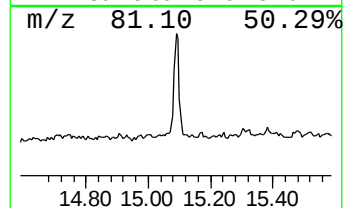
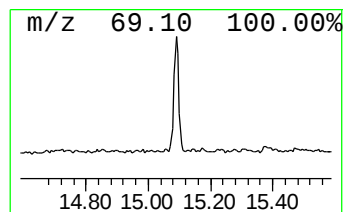
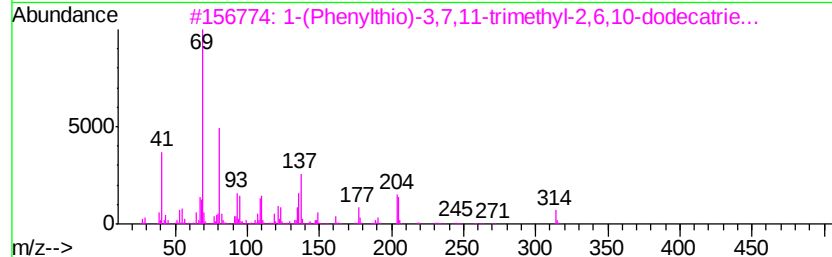
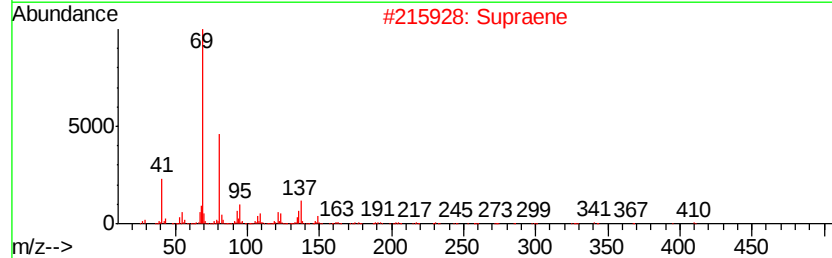
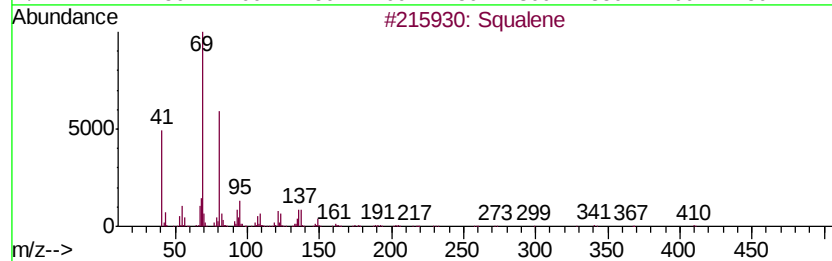
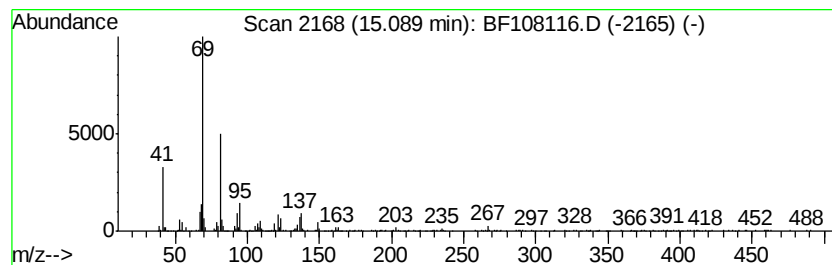
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Peak Number 4 Squalene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.09	5.75 ng	370305	Perylene-d12	15.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	94
2		Supraene	410	C30H50	007683-64-9	93
3		1-(Phenylthio)-3,7,11-trimethyl-...	314	C21H30S	028413-57-2	64
4		1,5-Heptadiene, 3,3,5-trimethyl-	138	C10H18	074630-29-8	64
5		6,11-Dimethyl-2,6,10-dodecatrien...	208	C14H24O	1000196-53-3	64



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.25	8.2	ng	647982	1	7.02	1590690	20.0
unknown6.79	6.79	41.6	ng	3305120	1	7.02	1590690	20.0
Squalene	15.09	5.8	ng	370305	6	15.81	1288600	20.0