

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
 Data File : BF108097.D
 Acq On : 10 Aug 2018 21:18
 Operator : JU/SJ
 Sample : J4272-19 2X
 Misc :
 ALS Vial : 21 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	5.248	491	495	505	rVB	641831	701125	18.52%	1.838%
2	5.637	557	561	565	rBV	3685369	3512164	92.78%	9.205%
3	6.631	725	730	733	rBV	3996338	3785537	100.00%	9.922%
4	6.660	734	735	741	rVB	61256	49350	1.30%	0.129%
5	6.784	752	756	760	rBV	3790638	3659996	96.68%	9.593%
6	7.019	792	796	800	rBV	2238053	1751877	46.28%	4.592%
7	7.178	819	823	826	rBV	3155952	2694398	71.18%	7.062%
8	7.578	886	891	894	rBV	2720931	2257274	59.63%	5.916%
9	8.301	1010	1014	1018	rBV	2410698	2088978	55.18%	5.475%
10	9.377	1192	1197	1200	rBV	4258747	3417762	90.28%	8.958%
11	9.766	1260	1263	1268	rBV	288652	225472	5.96%	0.591%
12	10.066	1310	1314	1318	rBV	2086040	1949052	51.49%	5.109%
13	10.101	1318	1320	1328	rVB	43175	43075	1.14%	0.113%
14	10.477	1381	1384	1387	rVB2	51392	43851	1.16%	0.115%
15	10.860	1445	1449	1459	rBV	1996030	1808656	47.78%	4.741%
16	10.954	1462	1465	1476	rVB	50928	95703	2.53%	0.251%
17	11.407	1539	1542	1544	rBV	51786	46305	1.22%	0.121%
18	11.566	1565	1569	1571	rBV	2134995	1770846	46.78%	4.641%
19	11.589	1571	1573	1576	rVB	304996	235437	6.22%	0.617%
20	11.642	1579	1582	1584	rVB	98803	83258	2.20%	0.218%
21	12.066	1651	1654	1657	rBV	136131	126570	3.34%	0.332%
22	12.095	1657	1659	1663	rVB2	72863	63684	1.68%	0.167%
23	12.183	1671	1674	1676	rBV2	87963	95463	2.52%	0.250%
24	12.624	1746	1749	1751	rBV3	72246	86288	2.28%	0.226%
25	12.707	1761	1763	1766	rBV2	38433	42000	1.11%	0.110%
26	12.783	1773	1776	1780	rBV	650025	570022	15.06%	1.494%
27	13.018	1810	1816	1821	rBV	596232	668175	17.65%	1.751%
28	13.154	1835	1839	1842	rBV	3115445	2671458	70.57%	7.002%
29	13.342	1869	1871	1873	rVB	76182	63046	1.67%	0.165%
30	14.224	2016	2021	2023	rBV2	1713685	1872621	49.47%	4.908%
31	14.248	2023	2025	2030	rVB	257409	229987	6.08%	0.603%
32	15.083	2165	2167	2171	rVB	311136	303993	8.03%	0.797%
33	15.795	2283	2288	2298	rVB	777307	1139591	30.10%	2.987%

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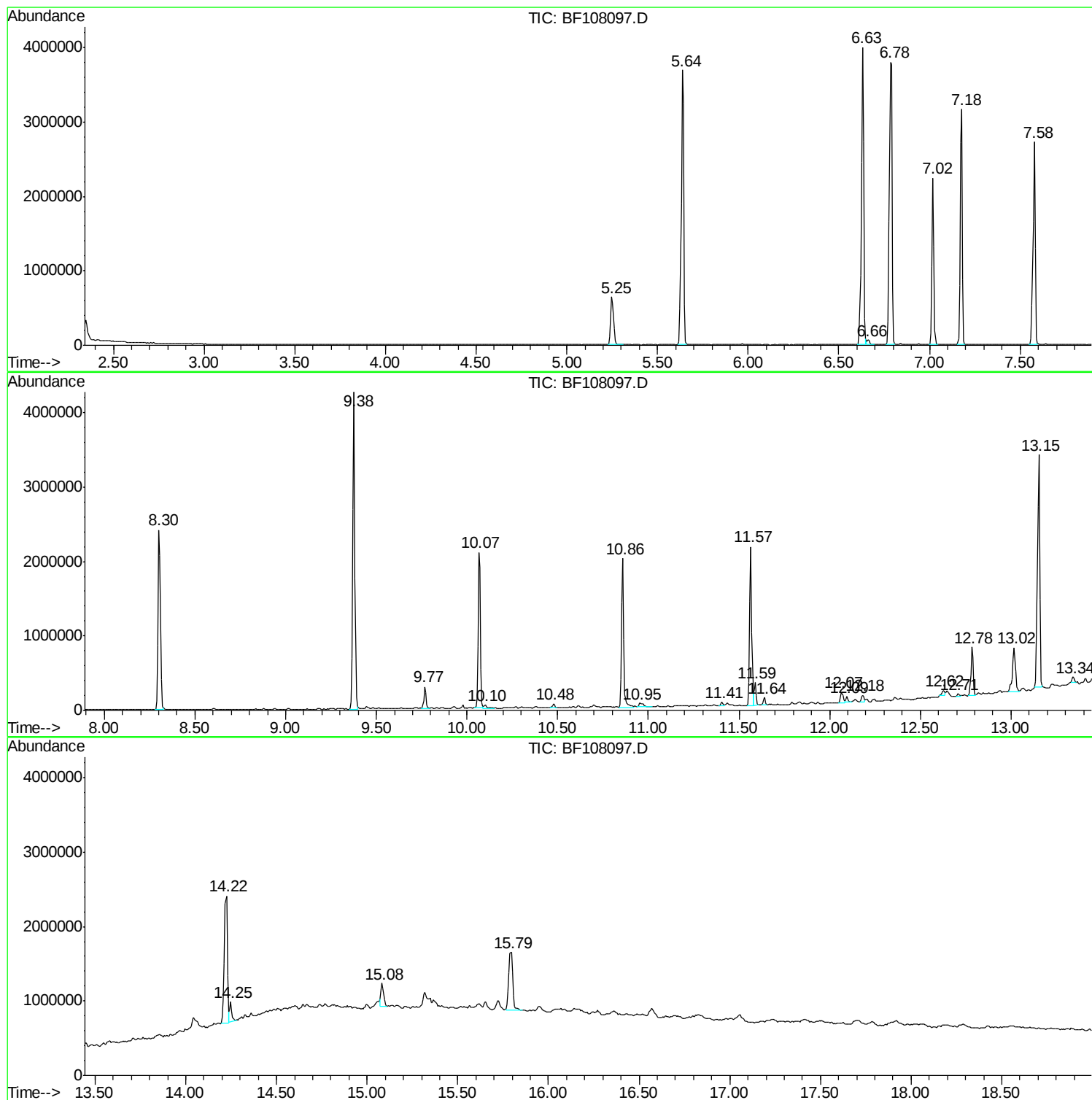
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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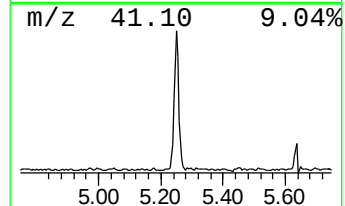
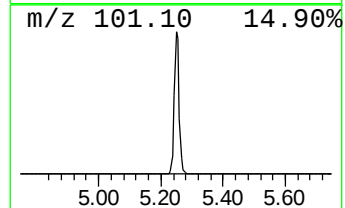
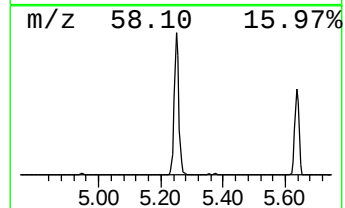
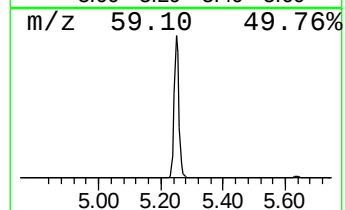
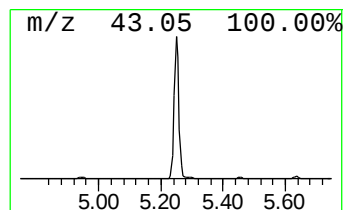
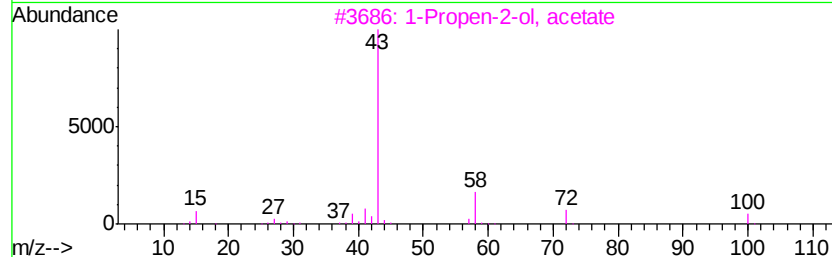
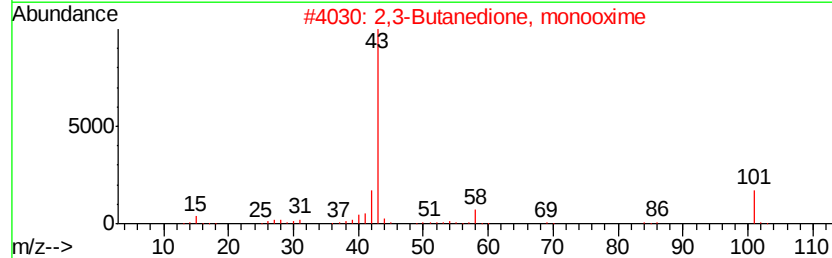
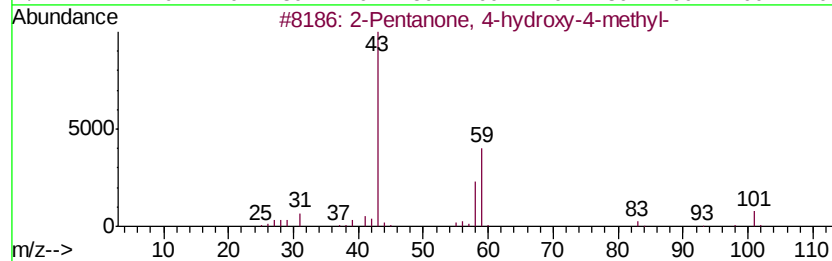
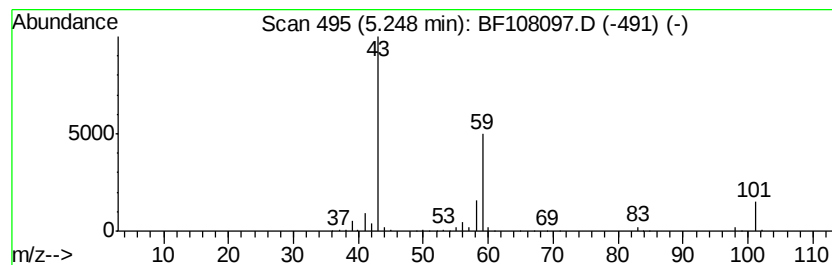
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.25	8.00 ng	701125	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	50
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4			Allyl acetate	100	C5H8O2	000591-87-7	9
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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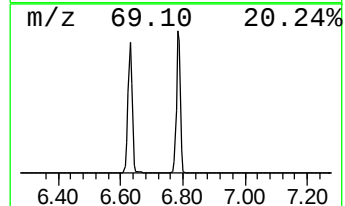
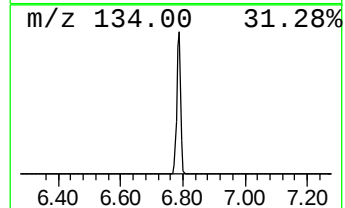
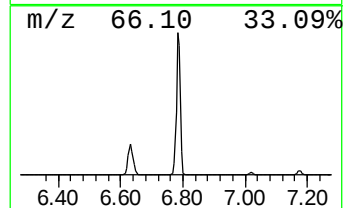
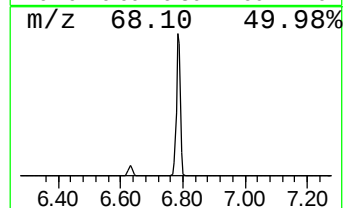
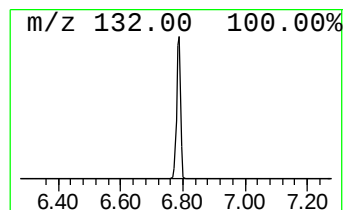
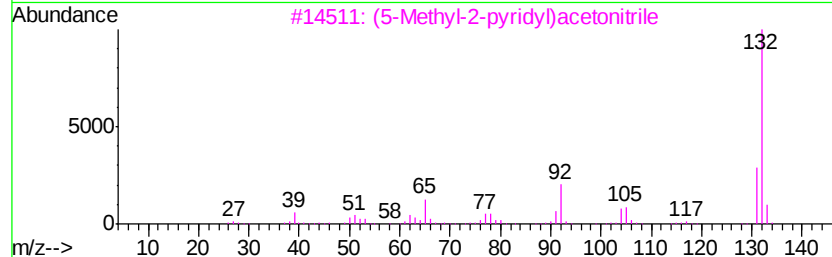
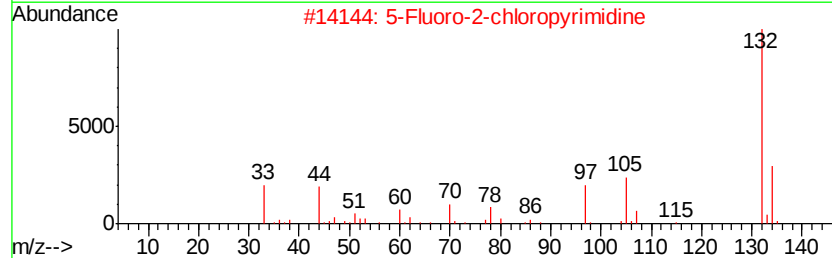
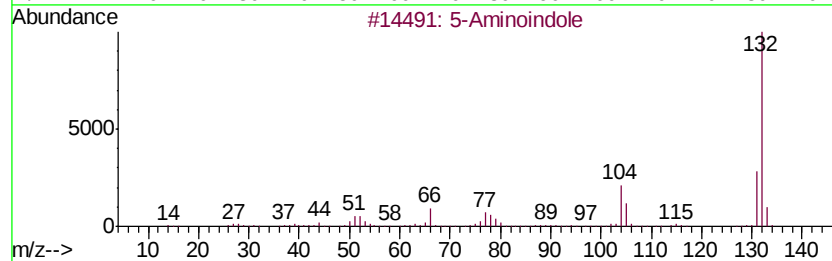
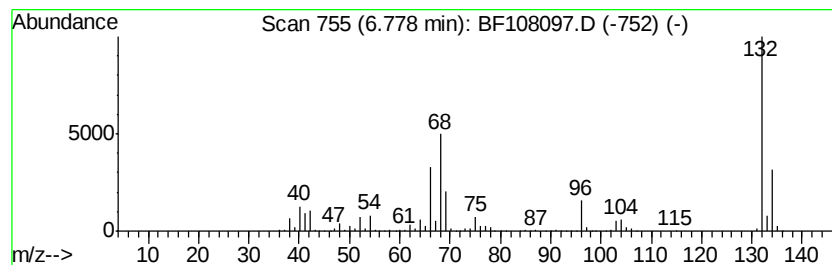
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Peak Number 2 unknown6.78 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.78	41.78 ng	3660000	1,4-Dichlorobenzene-d4	7.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		1,2,3-Trifluorobenzene	132	C6H3F3	001489-53-8	9



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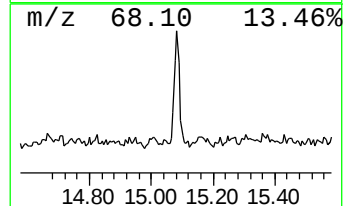
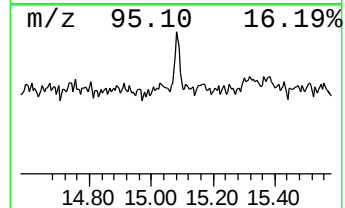
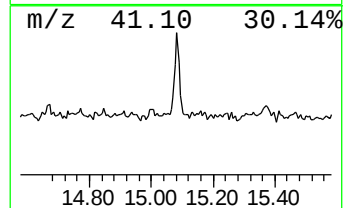
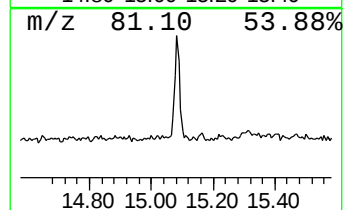
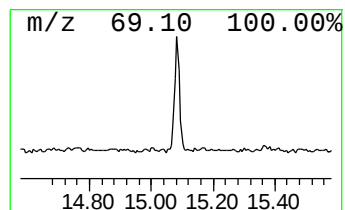
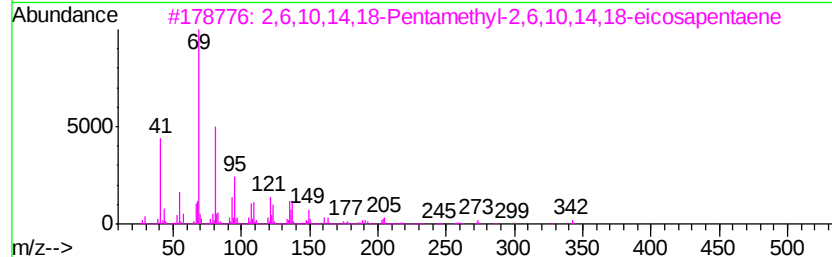
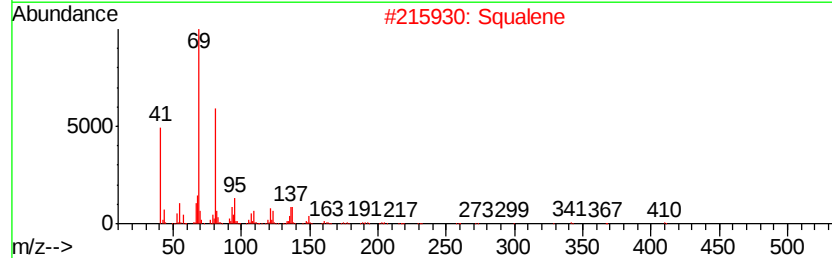
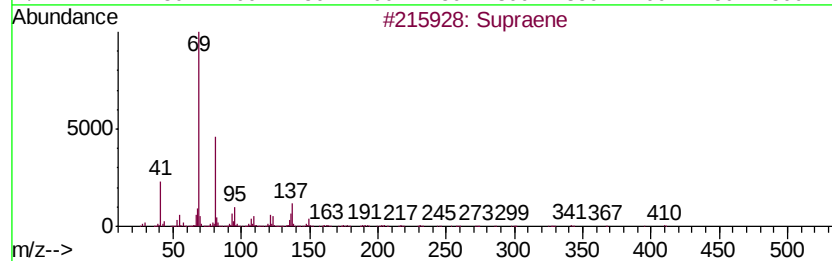
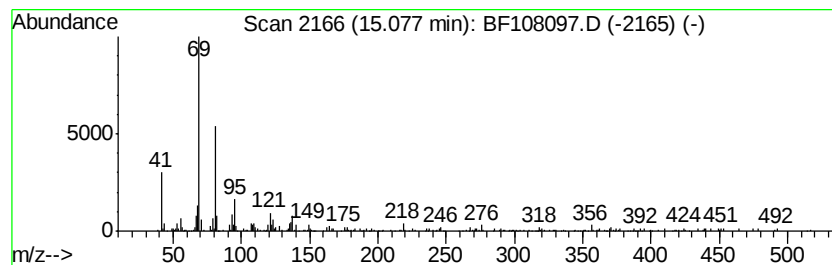
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Peak Number 4 Supraene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.08	5.34 ng	303993	Perylene-d12	15.79
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Supraene	410 C30H50	007683-64-9	93
2	Squalene	410 C30H50	000111-02-4	87
3	2,6,10,14,18-Pentamethyl-2,6,10,...	342 C25H42	075581-03-2	83
4	Farnesol isomer a	222 C15H26O	1000108-92-4	72
5	2,6,10-Dodecatrien-1-ol, 3,7,11-...	264 C17H28O2	004128-17-0	64



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.25	8.0	ng	701125	1	7.02	1751880	20.0
unknown6.78	6.78	41.8	ng	3660000	1	7.02	1751880	20.0
Supraene	15.08	5.3	ng	303993	6	15.79	1139590	20.0