

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082719\
 Data File : BG042521.D
 Acq On : 28 Aug 2019 1:56
 Operator : HP/JU
 Sample : K4516-17
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 T2-S-B1-B4

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_G\METHODS\8270-BG082219.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	3.117	3	5	26	rVB	897809	1243406	41.37%	7.700%
2	5.555	413	420	435	rBV	537819	872543	29.03%	5.403%
3	7.165	685	694	710	rBV	535816	963219	32.05%	5.965%
4	7.529	748	756	770	rBV	574844	1008613	33.56%	6.246%
5	7.999	828	836	843	rBV	118369	201485	6.70%	1.248%
6	8.323	884	891	900	rBV	470614	823657	27.41%	5.100%
7	9.169	1027	1035	1047	rBV	310089	590878	19.66%	3.659%
8	10.820	1310	1316	1330	rBV	137747	265186	8.82%	1.642%
9	13.276	1727	1734	1757	rBV	966716	1560775	51.94%	9.665%
10	14.104	1869	1875	1887	rBV	119517	193087	6.43%	1.196%
11	14.656	1962	1969	1981	rBV	266414	430283	14.32%	2.665%
12	16.149	2216	2223	2247	rBV	969347	1599874	53.24%	9.907%
13	16.360	2252	2259	2270	rVB2	13684	35184	1.17%	0.218%
14	17.412	2430	2438	2451	rBV2	380909	619432	20.61%	3.836%
15	17.529	2451	2458	2465	rVB4	19114	32299	1.07%	0.200%
16	19.468	2782	2788	2793	rBV	24507	38711	1.29%	0.240%
17	19.827	2838	2849	2854	rBV	25511	55659	1.85%	0.345%
18	20.021	2876	2882	2890	rBV	2208398	3005242	100.00%	18.610%
19	21.331	3101	3105	3109	rBV3	25311	46629	1.55%	0.289%
20	21.683	3158	3165	3179	rVB2	518574	896693	29.84%	5.553%
21	22.488	3298	3302	3308	rBV3	30715	45216	1.50%	0.280%
22	23.187	3416	3421	3428	rVB2	46932	88727	2.95%	0.549%
23	23.381	3447	3454	3464	rVB	184937	373932	12.44%	2.316%
24	23.998	3555	3559	3568	rVB2	41520	87400	2.91%	0.541%
25	24.927	3706	3717	3733	rBV2	323582	1007518	33.53%	6.239%
26	26.096	3910	3916	3925	rVB5	21392	62953	2.09%	0.390%

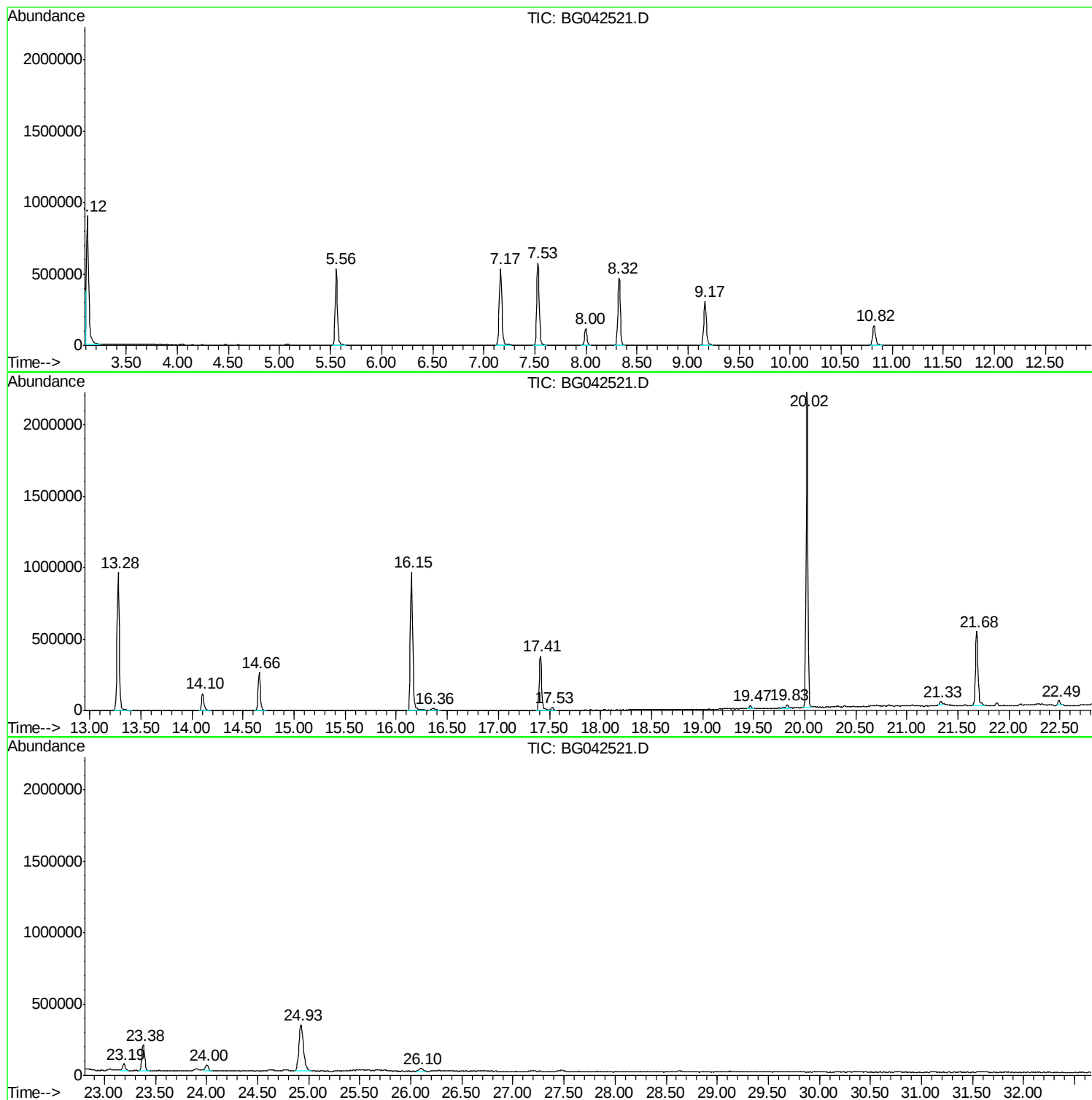
Sum of corrected areas: 16148601

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



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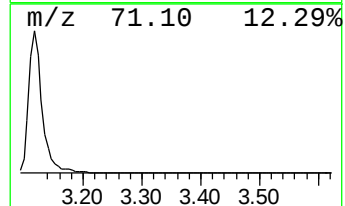
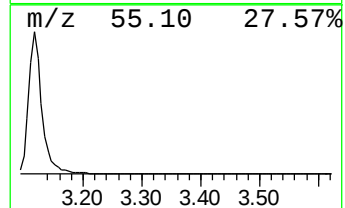
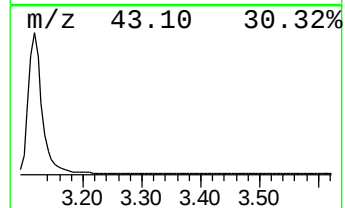
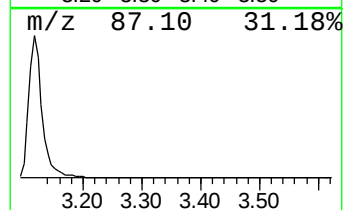
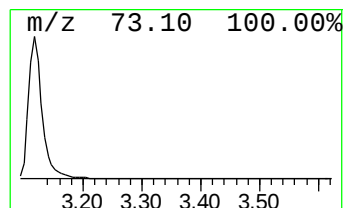
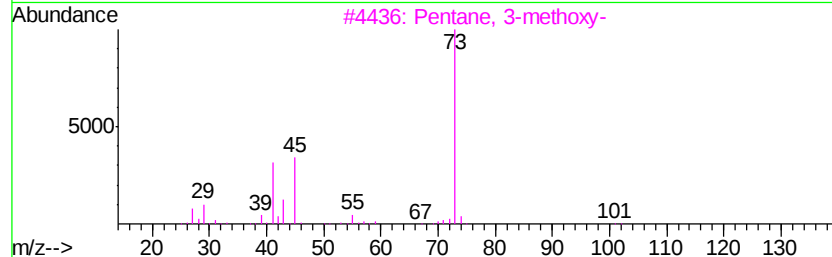
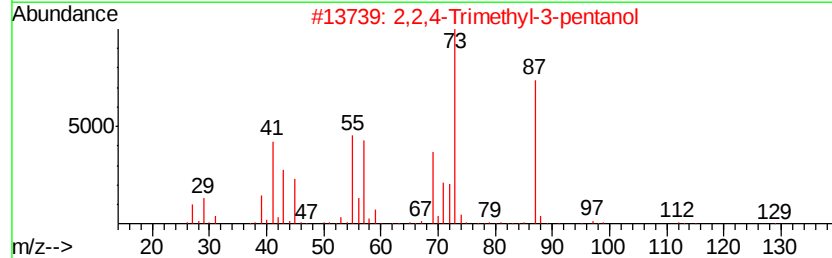
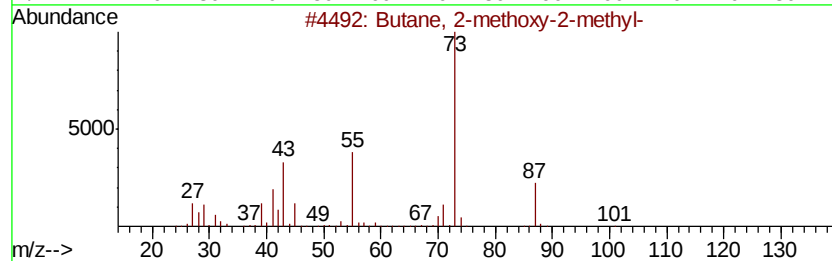
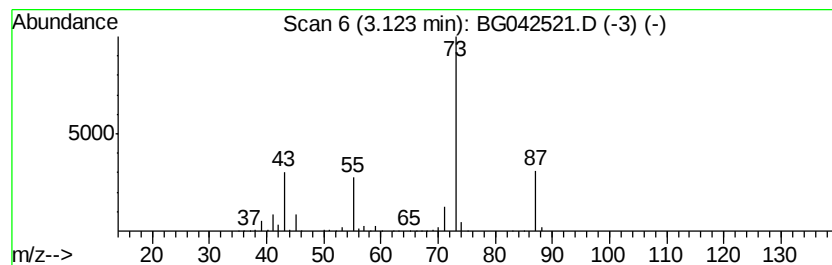
Instrument :
BNA_G
ClientSampleId :
T2-S-B1-B4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
3.12	123.42 ng	1243410	1,4-Dichlorobenzene-d4	8.00		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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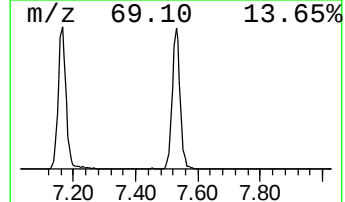
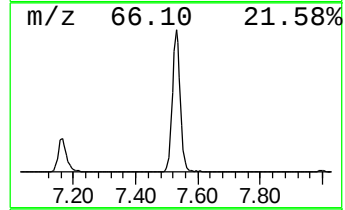
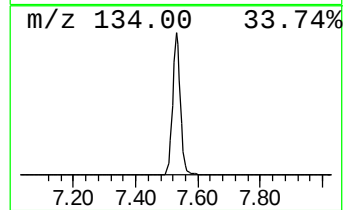
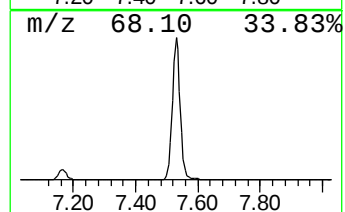
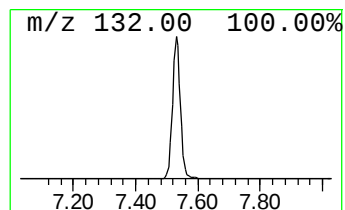
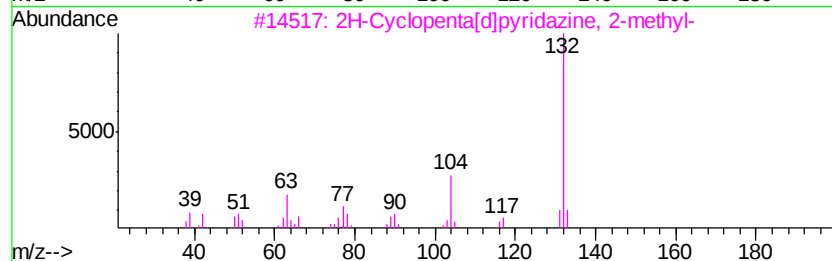
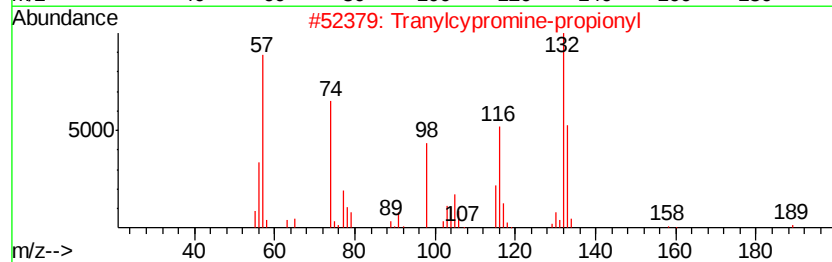
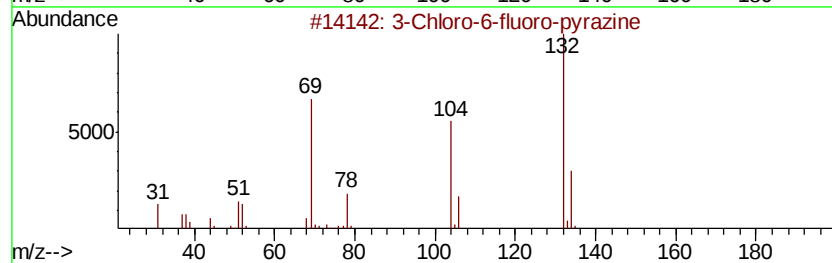
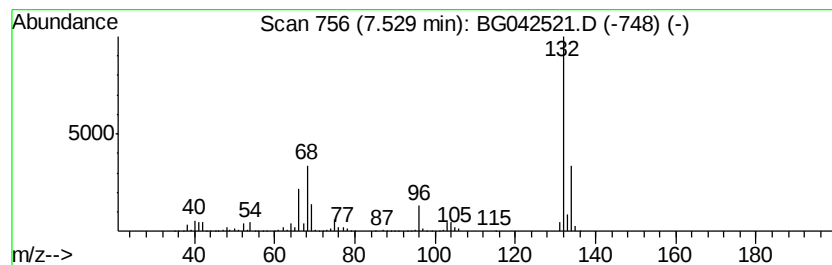
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Peak Number 2 unknown7.53 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.53	100.12 ng	1008610	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	16
3		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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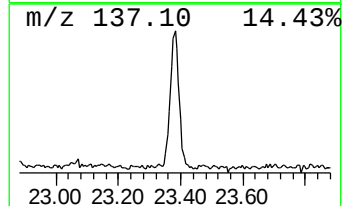
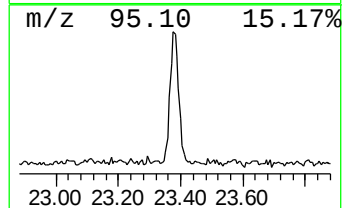
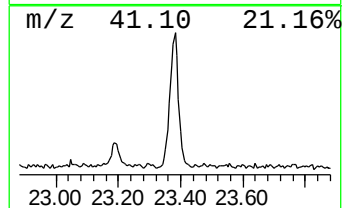
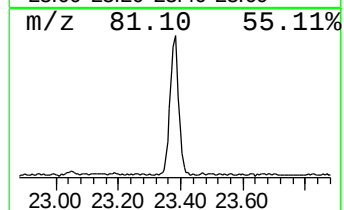
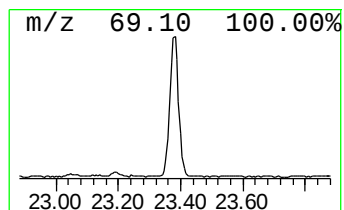
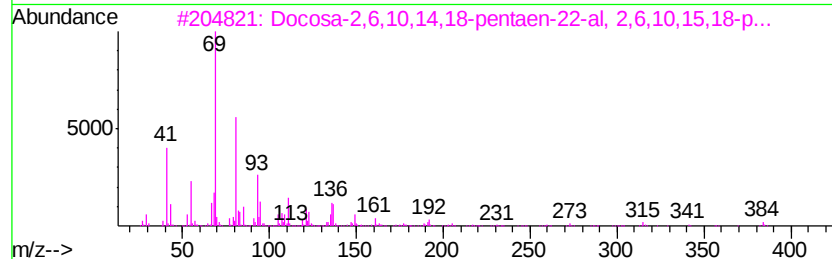
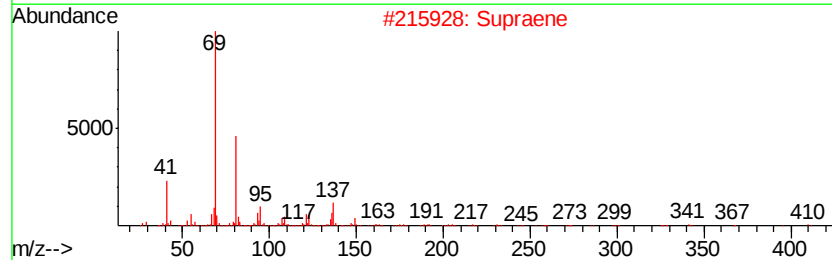
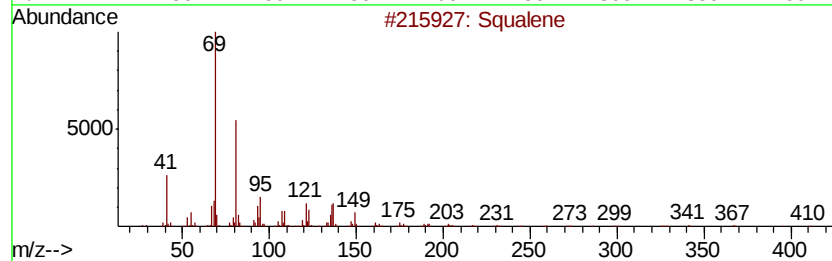
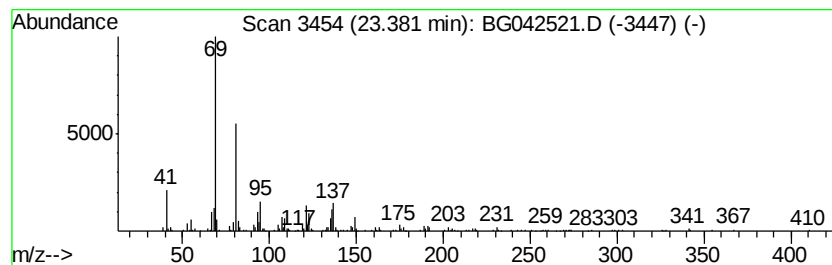
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 Peak Number 4 Squalene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.38	7.42 ng	373932	Perylene-d12	24.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	99
2		Supraene	410	C30H50	007683-64-9	98
3		Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	83
4		1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	74
5		1,5,9-Decatriene, 2,3,5,8-tetram...	192	C14H24	230646-72-7	72



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.12	123.4	ng	1243410	1	8.00	201485	20.0
unknown7.53	7.53	100.1	ng	1008610	1	8.00	201485	20.0
Squalene	23.38	7.4	ng	373932	6	24.93	1007520	20.0