

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030915\
 Data File : BF077648.D
 Acq On : 9 Mar 2015 20:01
 Operator : TP/IZ
 Sample : G1453-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SEP2

Quant Time: Mar 10 02:25:02 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 10 00:35:08 2015
 Response via : Initial Calibration

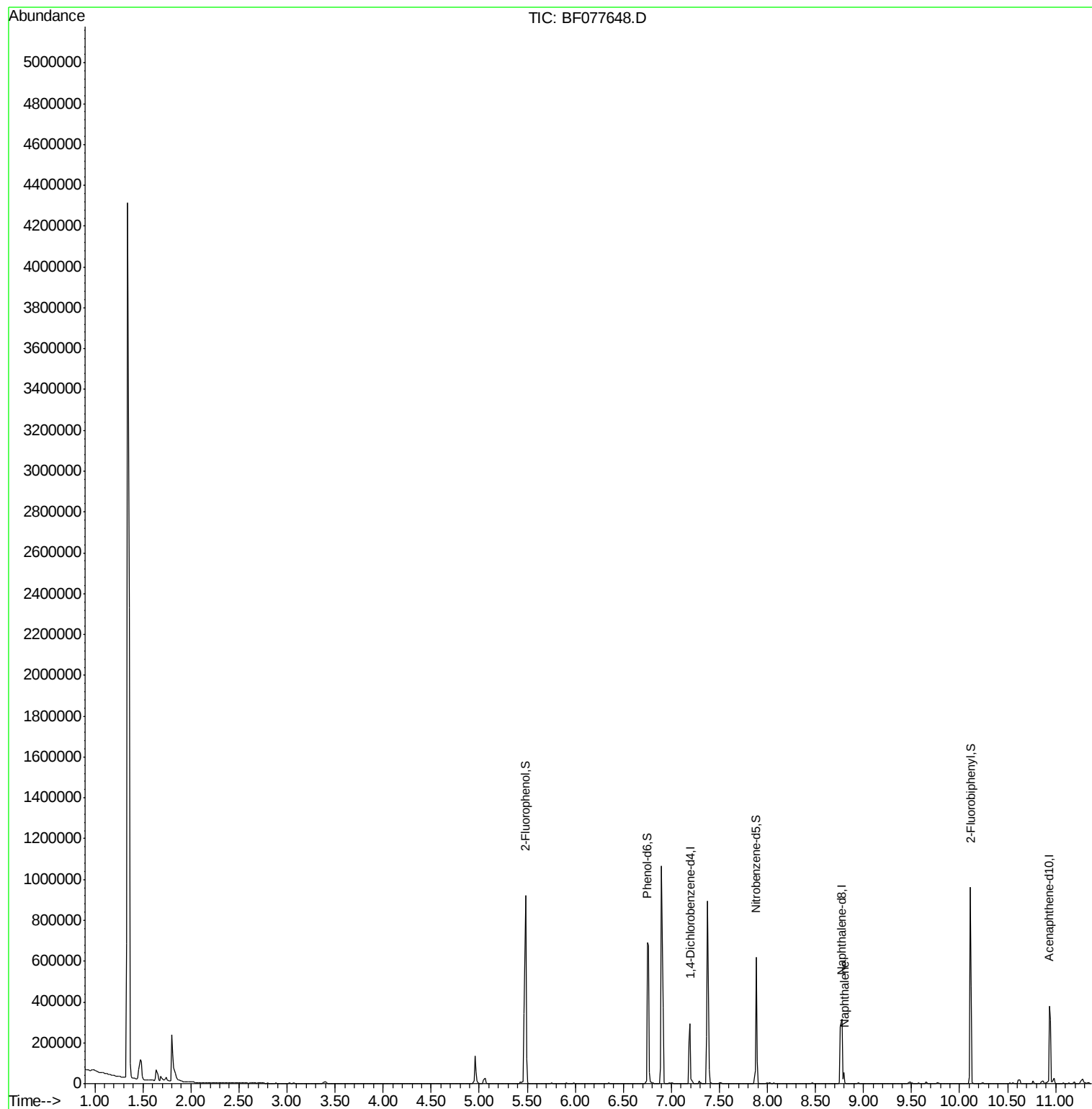
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.19	152	59364	20.00	ng	0.00
21) Naphthalene-d8	8.77	136	213179	20.00	ng	0.00
38) Acenaphthene-d10	10.94	164	107962	20.00	ng	0.00
63) Phenanthrene-d10	12.78	188	214481	20.00	ng	0.00
75) Chrysene-d12	16.07	240	243307	20.00	ng	0.01
86) Perylene-d12	17.87	264	262968	20.00	ng	0.13
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	293553	82.55	ng	0.00
7) Phenol-d6	6.76	99	374106	81.82	ng	0.00
23) Nitrobenzene-d5	7.88	82	190325	55.52	ng	0.00
41) 2,4,6-Tribromophenol	11.92	330	86576	83.28	ng	0.00
44) 2-Fluorobiphenyl	10.11	172	368192	53.18	ng	0.00
78) Terphenyl-d14	14.77	244	435471	41.84	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.79	128	26385	2.47	ng	99
70) Phenanthrene	12.81	178	145195	11.68	ng	98
71) Anthracene	12.87	178	36078	2.92	ng	97
74) Fluoranthene	14.28	202	243635	17.94	ng	98
77) Pyrene	14.56	202	223657	15.03	ng	97
80) Benzo(a)anthracene	16.06	228	110928	7.78	ng	98
82) Chrysene	16.10	228	102562	7.66	ng	97
85) Indeno(1,2,3-cd)pyrene	19.28	276	64442m	4.22	ng	
87) Benzo(b)fluoranthene	17.40	252	150260m	9.83	ng	
88) Benzo(k)fluoranthene	17.44	252	32821m	2.19	ng	
89) Benzo(a)pyrene	17.80	252	103453	7.10	ng	# 97
91) Benzo(a,h,i)perylene	19.70	276	64120	4.57	ng	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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