

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF101217\
 Data File : BF099415.D
 Acq On : 13 Oct 2017 00:43
 Operator : SJ/JU
 Sample : I5729-13
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 7E-(1-2)

Manual Integrations
 APPROVED

Quant Time: Oct 13 04:49:57 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 16:25:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	110102	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	430316	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	166834	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	230793	20.00	ng	0.00
75) Chrysene-d12	14.02	240	189642	20.00	ng	0.00
86) Perylene-d12	15.50	264	162702	20.00	ng	0.02

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System Monitoring Compounds

5) 2-Fluorophenol	5.48	112	828806	119.83	ng	0.01
7) Phenol-d6	6.49	99	982188	115.12	ng	0.00
23) Nitrobenzene-d5	7.42	82	591141	88.10	ng	0.00
41) 2,4,6-Tribromophenol	10.68	330	133165	97.55	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	961193	96.18	ng	0.00
78) Terphenyl-d14	12.96	244	624245	74.86	ng	0.00

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Target Compounds

						Qvalue
10) Phenol	6.50	94	27561	2.888	ng	# 70
31) Naphthalene	8.16	128	53083	2.627	ng	100
37) 2-Methylnaphthalene	8.85	142	29082	2.214	ng	98
49) Dimethylphthalate	9.60	163	156796	12.452	ng	98
51) Acenaphthene	9.92	154	45220	4.332	ng	98
54) Dibenzofuran	10.09	168	59988	4.316	ng	97
57) Fluorene	10.43	166	54407	4.870	ng	99
70) Phenanthrene	11.40	178	964704	78.090	ng	99
71) Anthracene	11.45	178	188374	14.903	ng	98
72) Carbazole	11.61	167	111817	9.033	ng	98
74) Fluoranthene	12.60	202	1483704	118.606	ng	100
77) Pyrene	12.83	202	1301809	90.023	ng	100
80) Benzo(a)anthracene	14.02	228	622570	54.215	ng	99
82) Chrysene	14.05	228	603600	55.211	ng	98
85) Indeno(1,2,3-cd)pyrene	16.98	276	223660	25.574	ng	98
87) Benzo(b)fluoranthene	15.07	252	707807m	70.755	ng	
88) Benzo(k)fluoranthene	15.09	252	187948m	19.022	ng	
89) Benzo(a)pyrene	15.44	252	408443	44.480	ng	# 95
90) Dibenzo(a,h)anthracene	16.98	278	58654m	7.981	ng	
91) Benzo(a,h,i)perylene	17.43	276	217972	30.007	ng	95

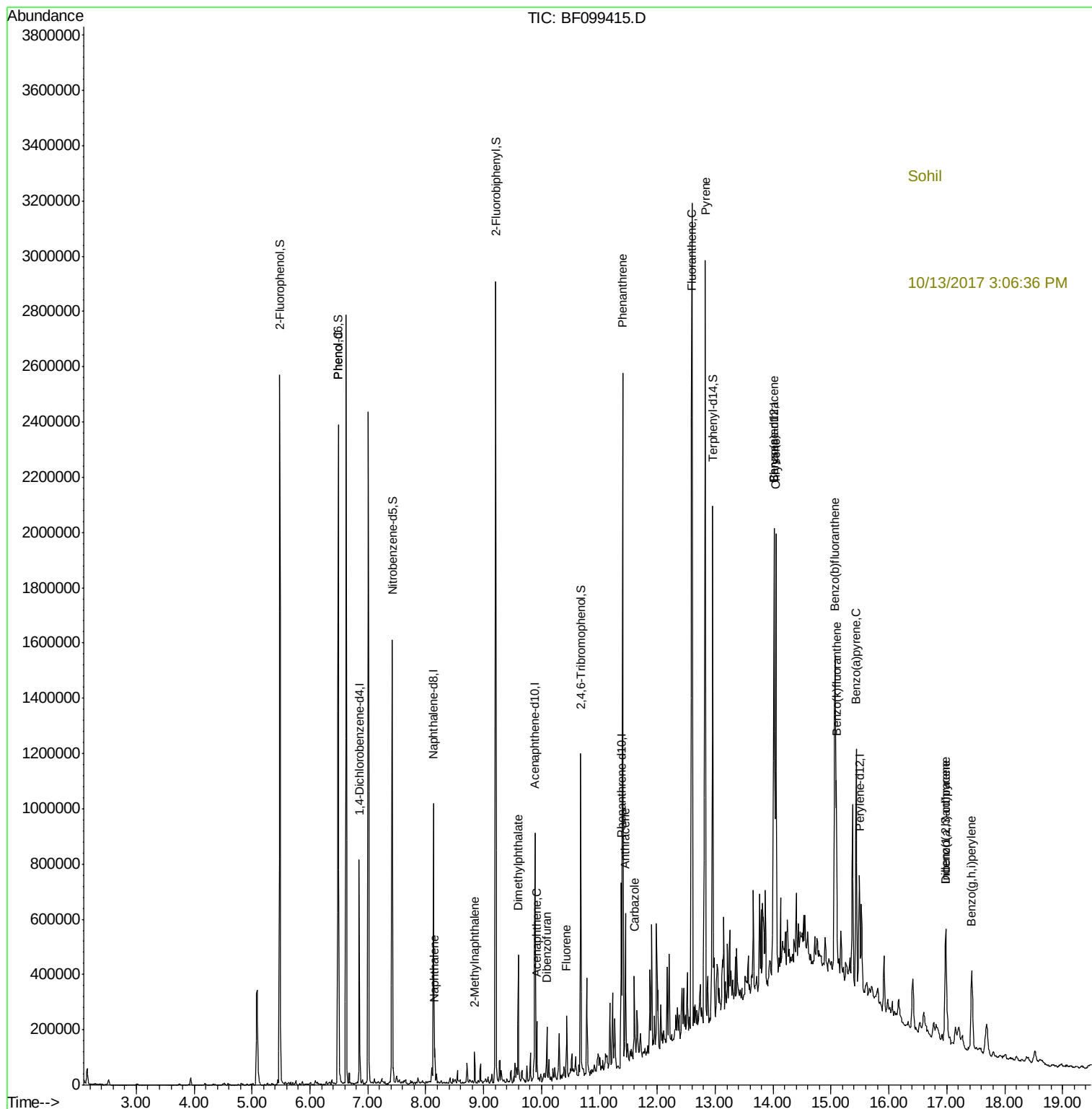
(#) = qualifier out of range (m) = manual integration (+) = signals summed

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