

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032117\
 Data File : BF093828.D
 Acq On : 22 Mar 2017 10:30
 Operator : SJ/MA
 Sample : I2318-09DL 50X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HP-PW-02DL

Quant Time: Mar 22 14:57:38 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 21 14:38:20 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	200883	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	805464	20.00	ng	0.00
38) Acenaphthene-d10	9.86	164	371882	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	568717	20.00	ng	0.00
75) Chrysene-d12	13.97	240	377286	20.00	ng	0.00
86) Perylene-d12	15.37	264	345650	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	27762	2.42	ng	-0.01
7) Phenol-d6	6.45	99	35966	2.53	ng	-0.01
23) Nitrobenzene-d5	7.38	82	19511	1.37	ng	-0.01
41) 2,4,6-Tribromophenol	10.64	330	5051	1.75	ng	-0.01
44) 2-Fluorobiphenyl	9.19	172	43768	1.98	ng	0.00
78) Terphenyl-d14	12.92	244	28519	1.54	ng	-0.01

Target Compounds				Qvalue		
31) Naphthalene	8.14	128	2748800	73.52	ng	99
37) 2-Methylnaphthalene	8.82	142	954651	39.12	ng	99
45) 1,1'-Biphenyl	9.28	154	142909	4.73	ng	95
48) Acenaphthylene	9.72	152	102090	2.71	ng	# 94
51) Acenaphthene	9.89	154	50311	2.23	ng	94
54) Dibenzofuran	10.06	168	72769	2.40	ng	# 87
57) Fluorene	10.40	166	248993	10.73	ng	99
70) Phenanthrene	11.36	178	946646	31.25	ng	99
71) Anthracene	11.42	178	258041	8.28	ng	96
74) Fluoranthene	12.55	202	477362	15.33	ng	96
77) Pyrene	12.77	202	563588	17.95	ng	98
80) Benzo(a)anthracene	13.96	228	172563m	7.30	ng	
82) Chrysene	13.99	228	153316m	6.49	ng	
85) Indeno(1,2,3-cd)pyrene	16.72	276	56543	2.92	ng	93
87) Benzo(b)fluoranthene	14.97	252	143639m	6.16	ng	
88) Benzo(k)fluoranthene	14.99	252	56509m	3.28	ng	
91) Benzo(g,h,i)perylene	17.12	276	53632	3.44	ng	95

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

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