

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120817\
 Data File : BF101253.D
 Acq On : 8 Dec 2017 23:07
 Operator : SJ/JU
 Sample : I6717-03
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB-3(0-10)

Quant Time: Dec 11 06:00:43 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 08 18:17:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	46934	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	178226	20.00	ng	0.00
38) Acenaphthene-d10	9.87	164	75536	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	132094	20.00	ng	0.00
75) Chrysene-d12	14.00	240	102379	20.00	ng	0.00
86) Perylene-d12	15.48	264	88541	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	312356	109.97	ng	0.01
7) Phenol-d6	6.47	99	376626	109.22	ng	0.00
23) Nitrobenzene-d5	7.39	82	218721	73.21	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	84361	92.97	ng	0.00
44) 2-Fluorobiphenyl	9.19	172	459296	83.19	ng	0.00
78) Terphenyl-d14	12.95	244	367386	78.49	ng	0.00
Target Compounds						
10) Phenol	6.47	94	9055	2.362	ng	# 50
49) Dimethylphthalate	9.58	163	55277	9.074	ng	# 99
70) Phenanthrene	11.38	178	51050	7.018	ng	96
74) Fluoranthene	12.57	202	68109	8.447	ng	96
77) Pyrene	12.80	202	61160	8.657	ng	98
80) Benzo(a)anthracene	13.99	228	29314	4.535	ng	99
82) Chrysene	14.03	228	26461	4.408	ng	99
85) Indeno(1,2,3-cd)pyrene	16.96	276	11780	2.207	ng	# 85
87) Benzo(b)fluoranthene	15.05	252	30147m	5.685	ng	
89) Benzo(a)pyrene	15.42	252	20901	4.335	ng	# 94
91) Benzo(g,h,i)perylene	17.40	276	11117	2.775	ng	94

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

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