

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042417\
 Data File : BF094584.D
 Acq On : 24 Apr 2017 14:44
 Operator : SJ/MA
 Sample : I2780-09
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PE-5

Quant Time: Apr 25 04:38:13 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 17 14:59:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.75	152	149381	20.00	ng	0.00
21) Naphthalene-d8	7.25	136	571696	20.00	ng	-0.01
38) Acenaphthene-d10	9.38	164	242374	20.00	ng	-0.02
63) Phenanthrene-d10	11.20	188	394893	20.00	ng	-0.02
75) Chrysene-d12	14.45	240	258033	20.00	ng	-0.02
86) Perylene-d12	16.08	264	237551	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.32	112	1216359	135.09	ng	0.01
7) Phenol-d6	5.40	99	1480430	139.47	ng	0.00
23) Nitrobenzene-d5	6.41	82	916412	98.98	ng	0.00
41) 2,4,6-Tribromophenol	10.36	330	260358	137.33	ng	-0.01
44) 2-Fluorobiphenyl	8.57	172	1406485	85.38	ng	-0.01
78) Terphenyl-d14	13.18	244	894559	92.67	ng	-0.01
Target Compounds						Qvalue
10) Phenol	5.40	94	93718	7.19	ng	95
49) Dimethylphthalate	9.07	163	332568	18.41	ng	98
74) Fluoranthene	12.69	202	64241	2.79	ng	97
77) Pyrene	12.97	202	58946	3.27	ng	95
80) Benzo(a)anthracene	14.44	228	31583	2.11	ng	99
82) Chrysene	14.48	228	32007	2.34	ng	98
87) Benzo(b)fluoranthene	15.67	252	52706m	3.63	ng	
89) Benzo(a)pyrene	16.02	252	33099	2.45	ng	98

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

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