

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042417\
 Data File : BF094581.D
 Acq On : 24 Apr 2017 13:21
 Operator : SJ/MA
 Sample : I2786-07
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BL-(3)

Quant Time: Apr 25 04:05:47 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	152075	20.00	ng	0.00
21) Naphthalene-d8	7.25	136	586564	20.00	ng	-0.01
38) Acenaphthene-d10	9.38	164	249157	20.00	ng	-0.02
63) Phenanthrene-d10	11.20	188	416047	20.00	ng	-0.02
75) Chrysene-d12	14.45	240	280566	20.00	ng	-0.02
86) Perylene-d12	16.08	264	246808	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.32	112	957695	104.48	ng	0.02
7) Phenol-d6	5.39	99	1188752	110.01	ng	0.00
23) Nitrobenzene-d5	6.41	82	737299	77.62	ng	0.00
41) 2,4,6-Tribromophenol	10.36	330	213306	109.45	ng	-0.01
44) 2-Fluorobiphenyl	8.57	172	1308741	77.28	ng	-0.01
78) Terphenyl-d14	13.18	244	930234	88.62	ng	-0.01
Target Compounds						Qvalue
10) Phenol	5.40	94	75270	5.67	ng	85
49) Dimethylphthalate	9.07	163	209398	11.28	ng	99
74) Fluoranthene	12.69	202	108210	4.46	ng	99
77) Pyrene	12.97	202	95214	4.86	ng	97
80) Benzo(a)anthracene	14.44	228	38459	2.36	ng	99
82) Chrysene	14.48	228	47934	3.22	ng	99
87) Benzo(b)fluoranthene	15.67	252	52540m	3.48	ng	
89) Benzo(a)pyrene	16.02	252	33708	2.40	ng	# 96

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

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