

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041917\
 Data File : BF094525.D
 Acq On : 19 Apr 2017 19:36
 Operator : SJ/MA
 Sample : I2750-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-U5-BASE

Quant Time: Apr 20 04:39:47 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	5.77	152	125503	20.00	ng	0.00
21) Naphthalene-d8	7.26	136	537131	20.00	ng	0.00
38) Acenaphthene-d10	9.40	164	211279	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	404891	20.00	ng	0.00
75) Chrysene-d12	14.46	240	279803	20.00	ng	0.00
86) Perylene-d12	16.08	264	238674	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.32	112	842972	111.44	ng	0.02
7) Phenol-d6	5.40	99	1084713	121.63	ng	0.00
23) Nitrobenzene-d5	6.42	82	683545	78.58	ng	0.00
41) 2,4,6-Tribromophenol	10.37	330	223659	135.34	ng	0.00
44) 2-Fluorobiphenyl	8.58	172	1230914	85.72	ng	0.00
78) Terphenyl-d14	13.19	244	970285	92.69	ng	0.00
Target Compounds						
10) Phenol	5.41	94	25973	2.37	ng	97
49) Dimethylphthalate	9.08	163	362769	23.04	ng	99
70) Phenanthrene	11.24	178	51953	2.28	ng	98
74) Fluoranthene	12.71	202	75157	3.18	ng	99
77) Pyrene	12.98	202	73149	3.74	ng	97
82) Chrysene	14.49	228	30648	2.06	ng	97
87) Benzo(b)fluoranthene	15.68	252	31065m	2.13	ng	

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

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