

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
 Data File : BF095328.D
 Acq On : 22 May 2017 22:42
 Operator : SJ/MA
 Sample : I3250-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W.P.RD-C.AVE(SP)

Quant Time: May 23 04:35:31 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 18 17:07:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	90241	20.00	ng	0.00
21) Naphthalene-d8	9.78	136	362593	20.00	ng	0.00
38) Acenaphthene-d10	12.61	164	142610	20.00	ng	-0.01
63) Phenanthrene-d10	15.01	188	220571	20.00	ng	-0.01
75) Chrysene-d12	18.68	240	204517	20.00	ng	-0.01
86) Perylene-d12	20.36	264	164010	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	649183	119.94	ng	-0.01
7) Phenol-d6	7.30	99	783731	120.68	ng	-0.01
23) Nitrobenzene-d5	8.66	82	438556	76.27	ng	-0.01
41) 2,4,6-Tribromophenol	13.91	330	123632	105.86	ng	-0.01
44) 2-Fluorobiphenyl	11.55	172	776133	78.33	ng	-0.01
78) Terphenyl-d14	17.33	244	498143	53.65	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	12.25	163	85642	7.32	ng	98
70) Phenanthrene	15.05	178	78381	6.18	ng	98
71) Anthracene	15.15	178	31212	2.41	ng	95
74) Fluoranthene	16.79	202	165232	12.71	ng	99
77) Pyrene	17.09	202	176999	11.57	ng	98
80) Benzo(a)anthracene	18.67	228	93367	7.84	ng	97
82) Chrysene	18.72	228	103057	8.95	ng	98
85) Indeno(1,2,3-cd)pyrene	21.71	276	36785	3.94	ng	93
87) Benzo(b)fluoranthene	19.95	252	107641m	10.62	ng	
88) Benzo(k)fluoranthene	19.98	252	41445m	4.24	ng	
89) Benzo(a)pyrene	20.31	252	78415	8.39	ng	98
91) Benzo(a,h,i)perylene	22.09	276	33878	4.47	ng	95

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

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