

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
 Data File : BF095322.D
 Acq On : 22 May 2017 19:30
 Operator : SJ/MA
 Sample : I3235-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TW1-TW2

Quant Time: May 23 04:06:26 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	80325	20.00	ng	0.00
21) Naphthalene-d8	9.78	136	332978	20.00	ng	0.00
38) Acenaphthene-d10	12.61	164	149374	20.00	ng	0.00
63) Phenanthrene-d10	15.01	188	214582	20.00	ng	-0.01
75) Chrysene-d12	18.68	240	189926	20.00	ng	-0.01
86) Perylene-d12	20.36	264	157807	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	558396	115.90	ng	-0.02
7) Phenol-d6	7.30	99	708520	122.57	ng	-0.02
23) Nitrobenzene-d5	8.66	82	382643	72.46	ng	-0.01
41) 2,4,6-Tribromophenol	13.91	330	123864	101.25	ng	-0.01
44) 2-Fluorobiphenyl	11.55	172	702167	67.66	ng	-0.01
78) Terphenyl-d14	17.32	244	374439	43.42	ng	-0.02
Target Compounds						
49) Dimethylphthalate	12.25	163	139075	11.34	ng	Qvalue 99
70) Phenanthrene	15.05	178	45009	3.65	ng	# 94
74) Fluoranthene	16.79	202	100371	7.94	ng	99
77) Pyrene	17.10	202	87228	6.14	ng	# 95
80) Benzo(a)anthracene	18.67	228	49959	4.52	ng	97
82) Chrysene	18.72	228	57932	5.42	ng	96
85) Indeno(1,2,3-cd)pyrene	21.71	276	22260	2.56	ng	95
87) Benzo(b)fluoranthene	19.96	252	65176m	6.68	ng	
88) Benzo(k)fluoranthene	19.98	252	28911m	3.07	ng	
89) Benzo(a)pyrene	20.31	252	43560	4.84	ng	95
91) Benzo(g,h,i)perylene	22.09	276	26781	3.67	ng	96

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

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