

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
 Data File : BF095326.D
 Acq On : 22 May 2017 21:38
 Operator : SJ/MA
 Sample : I3278-04
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RT-2669

Quant Time: May 23 14:17:12 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	83994	20.00	ng	0.00
21) Naphthalene-d8	9.78	136	353150	20.00	ng	0.00
38) Acenaphthene-d10	12.61	164	143191	20.00	ng	-0.01
63) Phenanthrene-d10	15.01	188	222427	20.00	ng	-0.01
75) Chrysene-d12	18.68	240	190499	20.00	ng	-0.01
86) Perylene-d12	20.36	264	158604	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.73	112	492075	97.67	ng	0.00
7) Phenol-d6	7.30	99	593334	98.16	ng	-0.01
23) Nitrobenzene-d5	8.66	82	341071	60.90	ng	-0.01
41) 2,4,6-Tribromophenol	13.91	330	101169	86.27	ng	0.00
44) 2-Fluorobiphenyl	11.55	172	656190	65.96	ng	-0.01
78) Terphenyl-d14	17.32	244	335128	38.75	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	12.25	163	77244	6.57	ng	100
70) Phenanthrene	15.05	178	60551	4.74	ng	96
74) Fluoranthene	16.79	202	137220	10.47	ng	96
77) Pyrene	17.09	202	120348	8.45	ng	# 94
80) Benzo(a)anthracene	18.67	228	49488	4.46	ng	99
82) Chrysene	18.72	228	66386m	6.19	ng	
85) Indeno(1,2,3-cd)pyrene	21.71	276	21380	2.46	ng	96
87) Benzo(b)fluoranthene	19.95	252	66497m	6.79	ng	
88) Benzo(k)fluoranthene	19.98	252	24725m	2.62	ng	
89) Benzo(a)pyrene	20.31	252	41586	4.60	ng	97
91) Benzo(g,h,i)perylene	22.11	276	20917	2.85	ng	94

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

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