

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042417\
 Data File : BF094586.D
 Acq On : 24 Apr 2017 15:39
 Operator : SJ/MA
 Sample : I2780-05 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PE-3

Quant Time: Apr 25 05:10:31 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	148321	20.00	ng	-0.01
21) Naphthalene-d8	7.24	136	565958	20.00	ng	-0.02
38) Acenaphthene-d10	9.38	164	238567	20.00	ng	-0.02
63) Phenanthrene-d10	11.20	188	359721	20.00	ng	-0.02
75) Chrysene-d12	14.45	240	266004	20.00	ng	-0.02
86) Perylene-d12	16.08	264	254386	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.30	112	191629	21.43	ng	0.00
7) Phenol-d6	5.37	99	239713	22.74	ng	-0.02
23) Nitrobenzene-d5	6.40	82	137952	15.05	ng	-0.02
41) 2,4,6-Tribromophenol	10.35	330	38070	20.40	ng	-0.02
44) 2-Fluorobiphenyl	8.57	172	273022	16.84	ng	-0.02
78) Terphenyl-d14	13.17	244	159181	16.00	ng	-0.03
Target Compounds						Qvalue
49) Dimethylphthalate	9.07	163	69736	3.92	ng	99
70) Phenanthrene	11.22	178	148241	7.32	ng	97
71) Anthracene	11.29	178	42562	2.03	ng	95
74) Fluoranthene	12.69	202	301091	14.34	ng	98
77) Pyrene	12.97	202	258459	13.91	ng	97
80) Benzo(a)anthracene	14.44	228	138450	8.95	ng	99
82) Chrysene	14.48	228	132335	9.37	ng	96
85) Indeno(1,2,3-cd)pyrene	17.35	276	52886	3.93	ng	98
87) Benzo(b)fluoranthene	15.68	252	160423m	10.31	ng	
88) Benzo(k)fluoranthene	15.70	252	50072m	3.40	ng	
89) Benzo(a)pyrene	16.02	252	109324	7.55	ng	# 96
91) Benzo(a,h,i)perylene	17.72	276	47858	3.91	ng	94

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

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