

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042617\
 Data File : BF094662.D
 Acq On : 26 Apr 2017 22:11
 Operator : SJ/MA
 Sample : I2824-04 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-W10-ESW

Quant Time: Apr 27 06:56:52 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	122562	20.00	ng	-0.01
21) Naphthalene-d8	7.24	136	486820	20.00	ng	-0.02
38) Acenaphthene-d10	9.38	164	193083	20.00	ng	-0.02
63) Phenanthrene-d10	11.20	188	311655	20.00	ng	-0.02
75) Chrysene-d12	14.45	240	254945	20.00	ng	-0.01
86) Perylene-d12	16.08	264	203883	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.30	112	150068	20.31	ng	0.00
7) Phenol-d6	5.37	99	184994	21.24	ng	-0.02
23) Nitrobenzene-d5	6.40	82	106928	13.56	ng	-0.02
41) 2,4,6-Tribromophenol	10.35	330	28403	18.81	ng	-0.02
44) 2-Fluorobiphenyl	8.57	172	220113	16.77	ng	-0.02
78) Terphenyl-d14	13.18	244	138916	14.56	ng	-0.02
Target Compounds						Qvalue
48) Acenaphthylene	9.21	152	54227	2.77	ng	98
70) Phenanthrene	11.23	178	134450	7.66	ng	97
71) Anthracene	11.30	178	57896	3.19	ng	97
74) Fluoranthene	12.69	202	380802	20.94	ng	98
77) Pyrene	12.97	202	335297	18.82	ng	97
80) Benzo(a)anthracene	14.44	228	156684m	10.57	ng	
82) Chrysene	14.48	228	171450	12.66	ng	96
83) Bis(2-ethylhexyl)phthalate	14.51	149	22007	2.10	ng	# 99
85) Indeno(1,2,3-cd)pyrene	17.36	276	72157	5.60	ng	95
87) Benzo(b)fluoranthene	15.68	252	218344m	17.51	ng	
88) Benzo(k)fluoranthene	15.71	252	74739m	6.33	ng	
89) Benzo(a)pyrene	16.02	252	110125	9.49	ng	# 94
90) Dibenzo(a,h)anthracene	17.37	278	20782	2.16	ng	# 86
91) Benzo(g,h,i)perylene	17.74	276	73009	7.44	ng	99

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

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