

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121416\
 Data File : BF091593.D
 Acq On : 14 Dec 2016 19:23
 Operator : UM/SJ
 Sample : H6006-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-13-5-10

Manual Integrations
 APPROVED

sohil
 12/15/2016 5:40:05 PM

Quant Time: Dec 15 01:17:01 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 09 19:00:21 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	297679	20.00	ng	-0.02
21) Naphthalene-d8	8.08	136	1167474	20.00	ng	-0.02
38) Acenaphthene-d10	9.84	164	616539	20.00	ng	-0.02
63) Phenanthrene-d10	11.31	188	945264	20.00	ng	-0.02
75) Chrysene-d12	13.94	240	614185	20.00	ng	-0.02
86) Perylene-d12	15.36	264	593078	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	1261616	71.40	ng	-0.02
7) Phenol-d6	6.43	99	1569702	71.26	ng	-0.02
23) Nitrobenzene-d5	7.36	82	1121305	53.61	ng	-0.02
41) 2,4,6-Tribromophenol	10.62	330	347983	67.95	ng	-0.01
44) 2-Fluorobiphenyl	9.16	172	1751986	49.10	ng	-0.02
78) Terphenyl-d14	12.90	244	1349544	49.86	ng	-0.02
Target Compounds						
10) Phenol	6.44	94	66696	2.59	ng	Qvalue 88
49) Dimethylphthalate	9.55	163	454399	11.64	ng	98
70) Phenanthrene	11.33	178	400138	8.55	ng	99
74) Fluoranthene	12.52	202	614578	12.54	ng	99
77) Pyrene	12.75	202	489404	10.04	ng	98
80) Benzo(a)anthracene	13.93	228	228266	6.40	ng	99
82) Chrysene	13.96	228	176036	4.72	ng	96
85) Indeno(1,2,3-cd)pyrene	16.71	276	119218	3.55	ng	97
87) Benzo(b)fluoranthene	14.96	252	251436m	7.11	ng	
88) Benzo(k)fluoranthene	14.98	252	83850m	2.70	ng	
89) Benzo(a)pyrene	15.29	252	170490	5.35	ng	# 95
91) Benzo(a,h,i)perylene	17.12	276	116437	4.01	ng	95

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

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