

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091216\
 Data File : BF090068.D
 Acq On : 12 Sep 2016 17:28
 Operator : UM/SJ
 Sample : H4744-04
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-1B

Quant Time: Sep 13 03:19:40 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF083116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 08 16:07:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.57	152	182736m	20.00	ng	-0.03
21) Naphthalene-d8	7.86	136	737755	20.00	ng	-0.03
38) Acenaphthene-d10	9.61	164	325576	20.00	ng	-0.02
63) Phenanthrene-d10	11.07	188	564577	20.00	ng	-0.03
75) Chrysene-d12	13.69	240	492103	20.00	ng	-0.03
86) Perylene-d12	15.03	264	353582	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.15	112	1219344	106.81	ng	-0.01
7) Phenol-d6	6.23	99	1378863	99.40	ng	-0.02
23) Nitrobenzene-d5	7.14	82	780446	61.35	ng	-0.03
41) 2,4,6-Tribromophenol	10.39	330	289228	90.04	ng	-0.03
44) 2-Fluorobiphenyl	8.95	172	1525351	76.91	ng	-0.02
78) Terphenyl-d14	12.66	244	1220640	61.27	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	9.34	163	508024	21.59	ng	99
70) Phenanthrene	11.10	178	104701	3.72	ng	96
74) Fluoranthene	12.27	202	187123	6.35	ng	96
77) Pyrene	12.50	202	170593	4.92	ng	99
80) Benzo(a)anthracene	13.68	228	75780m	2.51	ng	
82) Chrysene	13.71	228	86952m	3.09	ng	
87) Benzo(b)fluoranthene	14.67	252	73074m	3.30	ng	
89) Benzo(a)pyrene	14.97	252	53792	2.83	ng	# 93

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

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