

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121616\
 Data File : BF091682.D
 Acq On : 16 Dec 2016 19:37
 Operator : UM/SJ
 Sample : H6099-17
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUP-C-121316

Quant Time: Dec 16 23:26:20 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 16 00:35:41 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	155563	20.00	ng	-0.01
21) Naphthalene-d8	8.05	136	616632	20.00	ng	-0.02
38) Acenaphthene-d10	9.81	164	367575	20.00	ng	-0.01
63) Phenanthrene-d10	11.29	188	555043	20.00	ng	-0.02
75) Chrysene-d12	13.92	240	385934	20.00	ng	-0.02
86) Perylene-d12	15.32	264	318291	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	799279	86.56	ng	0.00
7) Phenol-d6	6.42	99	1063503	92.39	ng	-0.01
23) Nitrobenzene-d5	7.33	82	743927	67.33	ng	-0.02
41) 2,4,6-Tribromophenol	10.60	330	292569	95.83	ng	-0.01
44) 2-Fluorobiphenyl	9.14	172	1417257	66.62	ng	-0.01
78) Terphenyl-d14	12.88	244	1108082	65.15	ng	-0.02
Target Compounds						Qvalue
10) Phenol	6.43	94	51150	3.80	ng	77
49) Dimethylphthalate	9.53	163	134892	5.80	ng	# 98
70) Phenanthrene	11.31	178	78473	2.86	ng	96
74) Fluoranthene	12.50	202	113951	3.96	ng	96
77) Pyrene	12.73	202	93845	3.06	ng	99
80) Benzo(a)anthracene	13.91	228	53435	2.38	ng	97
87) Benzo(b)fluoranthene	14.92	252	57176m	3.01	ng	
89) Benzo(a)pyrene	15.25	252	43157	2.52	ng	98

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

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