

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121616\
 Data File : BF091683.D
 Acq On : 16 Dec 2016 20:04
 Operator : UM/SJ
 Sample : H6099-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP02-3FT

Quant Time: Dec 16 23:30:18 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	223087	20.00	ng	-0.01
21) Naphthalene-d8	8.05	136	856409	20.00	ng	-0.02
38) Acenaphthene-d10	9.81	164	491948	20.00	ng	-0.01
63) Phenanthrene-d10	11.29	188	813193	20.00	ng	-0.02
75) Chrysene-d12	13.92	240	666599	20.00	ng	-0.02
86) Perylene-d12	15.32	264	581658	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	1228818	92.80	ng	0.00
7) Phenol-d6	6.42	99	1868072	113.16	ng	-0.01
23) Nitrobenzene-d5	7.33	82	1048599	68.34	ng	-0.02
41) 2,4,6-Tribromophenol	10.60	330	474893	116.22	ng	-0.01
44) 2-Fluorobiphenyl	9.14	172	1762540	61.91	ng	-0.01
78) Terphenyl-d14	12.88	244	1606015	54.67	ng	-0.02
Target Compounds						Qvalue
10) Phenol	6.43	94	98406	5.10	ng	75
49) Dimethylphthalate	9.53	163	259778	8.34	ng	99
70) Phenanthrene	11.31	178	174933	4.34	ng	99
74) Fluoranthene	12.50	202	240556	5.70	ng	96
77) Pyrene	12.72	202	188980	3.57	ng	95
80) Benzo(a)anthracene	13.91	228	102173	2.64	ng	98
82) Chrysene	13.94	228	81463	2.01	ng	95
87) Benzo(b)fluoranthene	14.92	252	95170m	2.74	ng	

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

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