

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121616\
 Data File : BF091689.D
 Acq On : 16 Dec 2016 22:50
 Operator : UM/SJ
 Sample : H6099-16
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP13-1.5FT

Quant Time: Dec 19 13:02:02 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	168396	20.00	ng	-0.01
21) Naphthalene-d8	8.06	136	1014885	20.00	ng	-0.01
38) Acenaphthene-d10	9.81	164	497958	20.00	ng	-0.01
63) Phenanthrene-d10	11.29	188	640881	20.00	ng	-0.02
75) Chrysene-d12	13.92	240	375363	20.00	ng	-0.02
86) Perylene-d12	15.32	264	466169	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	1425195	142.59	ng	0.00
7) Phenol-d6	6.42	99	1066725	85.61	ng	-0.01
23) Nitrobenzene-d5	7.34	82	1282231	70.52	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	361779	87.47	ng	-0.01
44) 2-Fluorobiphenyl	9.14	172	1859774	64.53	ng	-0.01
78) Terphenyl-d14	12.88	244	1290252	78.00	ng	-0.02
Target Compounds						
10) Phenol	6.43	94	79441	5.45	ng	Qvalue 87
32) Benzoic acid	7.78	122	977	4.98	ng	# 80
49) Dimethylphthalate	9.53	163	199657	6.33	ng	97
74) Fluoranthene	12.50	202	90571	2.73	ng	98
77) Pyrene	12.73	202	76833	2.58	ng	97
87) Benzo(b)fluoranthene	14.92	252	73390m	2.64	ng	

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

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