

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110816\
 Data File : BF091111.D
 Acq On : 9 Nov 2016 5:33
 Operator : UM/SJ
 Sample : H5593-07
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TANK-GRAVEBOTTOM

Quant Time: Nov 09 06:24:14 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 12:59:29 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	185034	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	736184	20.00	ng	0.00
38) Acenaphthene-d10	9.70	164	343362	20.00	ng	0.00
63) Phenanthrene-d10	11.17	188	467389	20.00	ng	0.00
75) Chrysene-d12	13.80	240	333118	20.00	ng	-0.01
86) Perylene-d12	15.19	264	295028	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	1298194	115.87	ng	0.01
7) Phenol-d6	6.32	99	1663260	109.37	ng	0.00
23) Nitrobenzene-d5	7.23	82	989748	73.34	ng	-0.01
41) 2,4,6-Tribromophenol	10.49	330	294377	94.83	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	1473320	73.87	ng	-0.01
78) Terphenyl-d14	12.76	244	1118854	83.82	ng	0.00
Target Compounds						
10) Phenol	6.33	94	84685	5.03	ng	79
49) Dimethylphthalate	9.42	163	221086	9.82	ng	98

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

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