

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112916\
 Data File : BF091330.D
 Acq On : 29 Nov 2016 16:45
 Operator : UM/SJ
 Sample : H5696-01DL 2X
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GROUNDWATERDL

Quant Time: Nov 29 18:12:05 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 29 13:33:46 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	181951	20.00	ng	-0.02
21) Naphthalene-d8	8.15	136	768625	20.00	ng	-0.02
38) Acenaphthene-d10	9.90	164	390825	20.00	ng	-0.02
63) Phenanthrene-d10	11.38	188	632578	20.00	ng	-0.02
75) Chrysene-d12	14.01	240	355257	20.00	ng	-0.03
86) Perylene-d12	15.44	264	395556	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	145010	13.02	ng	0.00
7) Phenol-d6	6.48	99	128503	9.37	ng	-0.02
23) Nitrobenzene-d5	7.42	82	319960	23.33	ng	-0.03
41) 2,4,6-Tribromophenol	10.69	330	81907	22.14	ng	-0.02
44) 2-Fluorobiphenyl	9.22	172	654727	30.33	ng	-0.02
78) Terphenyl-d14	12.96	244	324570	19.86	ng	-0.03
Target Compounds						
10) Phenol	6.49	94	119100	7.38	ng	Qvalue 92
32) Benzoic acid	7.85	122	24042	2.72	ng	92
83) Bis(2-ethylhexyl)phthalate	14.01	149	570918	44.62	ng	# 94

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

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