

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111117\
 Data File : BG030955.D
 Acq On : 11 Nov 2017 22:45
 Operator : SJ/JU
 Sample : I6436-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 153RD-63RD-ST-DRUM

Quant Time: Nov 13 05:56:43 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 10 15:38:25 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	38157	20.00	ng	-0.01
21) Naphthalene-d8	10.95	136	173251	20.00	ng	-0.01
38) Acenaphthene-d10	14.76	164	129533	20.00	ng	-0.01
63) Phenanthrene-d10	17.50	188	335816	20.00	ng	-0.01
75) Chrysene-d12	21.77	240	408745	20.00	ng	-0.01
86) Perylene-d12	25.07	264	411844	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.71	112	151497	68.08	ng	0.00
7) Phenol-d6	7.30	99	121107	40.40	ng	0.00
23) Nitrobenzene-d5	9.31	82	288441	98.65	ng	-0.01
41) 2,4,6-Tribromophenol	16.25	330	328990	157.00	ng	0.00
44) 2-Fluorobiphenyl	13.38	172	897924	99.61	ng	-0.01
78) Terphenyl-d14	20.09	244	1646973	88.97	ng	0.00
Target Compounds						
10) Phenol	7.33	94	8765	2.815	ng	93
27) 2,4-Dimethylphenol	10.16	122	18906	8.180	ng	91
32) Benzoic acid	10.24	122	11805	11.071	ng	# 76
49) Dimethylphthalate	14.20	163	113902	10.170	ng	99

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

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