

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG010918\
 Data File : BG031972.D
 Acq On : 10 Jan 2018 8:18
 Operator : SJ/JU
 Sample : J1057-02
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OWS-2B

Quant Time: Jan 10 09:56:21 2018
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 05 15:31:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.78	152	65146	20.00	ng	-0.01
21) Naphthalene-d8	11.66	136	271954	20.00	ng	-0.01
38) Acenaphthene-d10	15.41	164	184321	20.00	ng	-0.02
63) Phenanthrene-d10	18.14	188	439816	20.00	ng	-0.02
75) Chrysene-d12	22.48	240	507568	20.00	ng	-0.04
86) Perylene-d12	26.20	264	513070	20.00	ng	-0.09
System Monitoring Compounds						
5) 2-Fluorophenol	6.21	112	407449	113.92	ng	0.00
7) Phenol-d6	7.89	99	468830	100.96	ng	0.00
23) Nitrobenzene-d5	9.97	82	351269	97.54	ng	-0.01
41) 2,4,6-Tribromophenol	16.88	330	381458	135.63	ng	-0.02
44) 2-Fluorobiphenyl	14.04	172	1245641	84.82	ng	-0.01
78) Terphenyl-d14	20.67	244	2018504	90.86	ng	-0.02
Target Compounds						
10) Phenol	7.91	94	34921	7.448	ng	95
20) 3+4-Methylphenols	9.54	107	54379	11.406	ng	94
49) Dimethylphthalate	14.83	163	40239	2.478	ng	# 95

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

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