

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG011718\
 Data File : BG032132.D
 Acq On : 18 Jan 2018 10:53
 Operator : SJ/JU
 Sample : J1158-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OWS-2

Quant Time: Jan 18 13:04:43 2018
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG011218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 18 10:13:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.75	152	43792	20.00	ng	0.00
21) Naphthalene-d8	11.63	136	171697	20.00	ng	0.00
38) Acenaphthene-d10	15.38	164	130459	20.00	ng	0.00
63) Phenanthrene-d10	18.11	188	350142	20.00	ng	0.00
75) Chrysene-d12	22.45	240	448104	20.00	ng	0.00
86) Perylene-d12	26.15	264	478251	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.17	112	299421	125.05	ng	0.00
7) Phenol-d6	7.85	99	344993	114.40	ng	0.00
23) Nitrobenzene-d5	9.94	82	256549	101.09	ng	0.00
41) 2,4,6-Tribromophenol	16.85	330	345306	159.95	ng	0.00
44) 2-Fluorobiphenyl	14.01	172	926134	88.02	ng	0.00
78) Terphenyl-d14	20.65	244	1632552	80.44	ng	0.00
Target Compounds						
10) Phenol	7.88	94	26071	8.776	ng	94
20) 3+4-Methylphenols	9.50	107	140598	44.706	ng	98
32) Benzoic acid	10.93	122	153974	77.063	ng	87

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

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