

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG012918\
 Data File : BG032430.D
 Acq On : 30 Jan 2018 00:58
 Operator : SJ/JU
 Sample : J1022-11
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SU-03-012618

Manual Integrations
 APPROVED

Sohil
 1/30/2018 6:23:07 PM

Quant Time: Jan 30 05:51:37 2018
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG012918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 29 15:32:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.72	152	28401	20.00	ng	0.00
21) Naphthalene-d8	11.60	136	118781	20.00	ng	0.00
38) Acenaphthene-d10	15.36	164	90636	20.00	ng	0.00
63) Phenanthrene-d10	18.09	188	249079	20.00	ng	0.00
75) Chrysene-d12	22.42	240	301811	20.00	ng	0.00
86) Perylene-d12	26.10	264	317930	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.16	112	216415	153.05	ng	0.00
7) Phenol-d6	7.84	99	249224	132.08	ng	0.00
23) Nitrobenzene-d5	9.92	82	197678	103.95	ng	0.00
41) 2,4,6-Tribromophenol	16.83	330	270190	156.61	ng	0.00
44) 2-Fluorobiphenyl	13.98	172	738201	96.80	ng	0.00
78) Terphenyl-d14	20.63	244	1496770	106.23	ng	0.00
Target Compounds						
31) Naphthalene	11.66	128	30129	5.193	ng	Qvalue 96
32) Benzoic acid	10.80	122	7618m	12.713	ng	
58) 2,3,4,6-Tetrachlorophenol	15.97	232	6390	2.410	ng	# 74
76) Benzidine	20.26	184	14573	2.459	ng	# 95

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

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