

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111117\
 Data File : BG030961.D
 Acq On : 12 Nov 2017 2:29
 Operator : SJ/JU
 Sample : I6301-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JC-02-110917-B

Manual Integrations
 APPROVED

SOHIL
 11/13/2017 4:55:36 PM

Quant Time: Nov 13 06:22:47 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	41961	20.00	ng	-0.01
21) Naphthalene-d8	10.96	136	173863	20.00	ng	-0.01
38) Acenaphthene-d10	14.76	164	124408	20.00	ng	-0.02
63) Phenanthrene-d10	17.50	188	358685	20.00	ng	-0.01
75) Chrysene-d12	21.77	240	466359	20.00	ng	-0.01
86) Perylene-d12	25.07	264	457516	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.70	112	336213	137.40	ng	0.00
7) Phenol-d6	7.31	99	392774	119.14	ng	0.00
23) Nitrobenzene-d5	9.31	82	296124	100.93	ng	-0.02
41) 2,4,6-Tribromophenol	16.24	330	355946	176.87	ng	-0.01
44) 2-Fluorobiphenyl	13.38	172	887696	102.53	ng	-0.02
78) Terphenyl-d14	20.09	244	1914654	90.65	ng	0.00
Target Compounds						
31) Naphthalene	11.01	128	42900	5.019	ng	98
32) Benzoic acid	10.28	122	837	6.044	ng	# 81
51) Acenaphthene	14.82	154	19534	2.567	ng	99
54) Dibenzofuran	15.16	168	24629	2.032	ng	97
57) Fluorene	15.80	166	20225	2.056	ng	# 96
69) Pentachlorophenol	17.17	266	11105m	4.162	ng	
70) Phenanthrene	17.54	178	50613	2.710	ng	99

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

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