

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071218\
 Data File : BG035610.D
 Acq On : 13 Jul 2018 1:04
 Operator : JU/SJ
 Sample : J3947-06
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PAR-3

Quant Time: Jul 13 04:52:43 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 12 14:43:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	43055	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	162073	20.00	ng	0.00
38) Acenaphthene-d10	14.67	164	119299	20.00	ng	0.00
63) Phenanthrene-d10	17.42	188	332850	20.00	ng	0.00
75) Chrysene-d12	21.68	240	367258	20.00	ng	0.00
86) Perylene-d12	24.90	264	335216	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	313562	120.91	ng	0.00
7) Phenol-d6	7.22	99	405899	104.91	ng	0.00
23) Nitrobenzene-d5	9.22	82	338824	87.33	ng	0.00
41) 2,4,6-Tribromophenol	16.17	330	221934	141.14	ng	0.00
44) 2-Fluorobiphenyl	13.30	172	745193	83.69	ng	0.00
78) Terphenyl-d14	20.03	244	1465842	87.98	ng	0.00
Target Compounds						
31) Naphthalene	10.91	128	28051	3.323	ng	Qvalue # 90
51) Acenaphthene	14.74	154	20248	2.697	ng	96
70) Phenanthrene	17.46	178	41089	2.474	ng	96

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

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