

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080316\
 Data File : BG023426.D
 Acq On : 3 Aug 2016 15:17
 Operator : UM/SJ
 Sample : H4287-12
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-61-7.0-8.0

Quant Time: Aug 04 03:27:09 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG080116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 19:22:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	107527	20.00	ng	0.00
21) Naphthalene-d8	10.95	136	473884	20.00	ng	0.00
38) Acenaphthene-d10	14.78	164	339353	20.00	ng	0.00
63) Phenanthrene-d10	17.54	188	890601	20.00	ng	0.00
75) Chrysene-d12	21.86	240	981529	20.00	ng	0.00
86) Perylene-d12	25.24	264	938780	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	962055	150.60	ng	0.00
7) Phenol-d6	7.27	99	1452001	145.94	ng	0.00
23) Nitrobenzene-d5	9.29	82	970970	101.86	ng	0.00
41) 2,4,6-Tribromophenol	16.28	330	721471	145.35	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	1853763	92.14	ng	0.00
78) Terphenyl-d14	20.15	244	2831972	91.83	ng	0.00
Target Compounds						Qvalue
10) Phenol	7.30	94	23245	2.18	ng	90
31) Naphthalene	11.00	128	286275	11.86	ng	95
37) 2-Methylnaphthalene	12.60	142	225099	12.27	ng	# 91
48) Acenaphthylene	14.50	152	256511	8.09	ng	98
49) Dimethylphthalate	14.22	163	1023263	39.30	ng	98
51) Acenaphthene	14.85	154	141758	7.05	ng	97
57) Fluorene	15.83	166	106435	4.18	ng	95

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

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