

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081516\
 Data File : BG023721.D
 Acq On : 15 Aug 2016 14:43
 Operator : UM/SJ
 Sample : H4452-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S-2

Quant Time: Aug 15 17:43:40 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	134513	20.00	ng	-0.02
21) Naphthalene-d8	10.94	136	585688	20.00	ng	0.00
38) Acenaphthene-d10	14.78	164	327128	20.00	ng	0.00
63) Phenanthrene-d10	17.55	188	644998	20.00	ng	0.00
75) Chrysene-d12	21.88	240	716272	20.00	ng	0.01
86) Perylene-d12	25.29	264	711319	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.65	112	1210614	151.50	ng	0.00
7) Phenol-d6	7.27	99	1724198	138.53	ng	0.00
23) Nitrobenzene-d5	9.29	82	1223117	103.82	ng	0.00
41) 2,4,6-Tribromophenol	16.28	330	602504	125.92	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	1991672	102.69	ng	0.00
78) Terphenyl-d14	20.16	244	2090327	93.17	ng	0.00
Target Compounds						
15) Benzyl Alcohol	8.43	79	21787	2.31	ng	# 75
49) Dimethylphthalate	14.22	163	756076	30.12	ng	99
80) Benzo(a)anthracene	21.86	228	87001	2.20	ng	# 89
82) Chrysene	21.92	228	75291	2.00	ng	94
87) Benzo(b)fluoranthene	24.20	252	106793	2.67	ng	# 98
89) Benzo(a)pyrene	25.12	252	78518	2.06	ng	# 91

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

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