

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071016\
 Data File : BG022986.D
 Acq On : 10 Jul 2016 5:29
 Operator : UM/SJ
 Sample : H3915-15
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P7-B4-0-6-C

Quant Time: Jul 11 03:37:37 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG070816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 07 16:51:07 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.15	152	97543	20.00	ng	0.00
21) Naphthalene-d8	10.98	136	452617	20.00	ng	0.00
38) Acenaphthene-d10	14.80	164	346675	20.00	ng	0.00
63) Phenanthrene-d10	17.55	188	822142	20.00	ng	0.00
75) Chrysene-d12	21.85	240	890896	20.00	ng	0.00
86) Perylene-d12	25.20	264	910841	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.70	112	446610	74.88	ng	0.00
7) Phenol-d6	7.31	99	747369	80.01	ng	0.00
23) Nitrobenzene-d5	9.32	82	557652	52.76	ng	0.00
41) 2,4,6-Tribromophenol	16.28	330	377324	60.56	ng	-0.01
44) 2-Fluorobiphenyl	13.42	172	1285927	58.43	ng	0.00
78) Terphenyl-d14	20.16	244	1900838	60.47	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.24	163	321192	11.34	ng	98
74) Fluoranthene	19.60	202	181951	3.68	ng	96
77) Pyrene	19.96	202	151555	3.03	ng	# 89
82) Chrysene	21.89	228	103210	2.21	ng	97
83) Bis(2-ethylhexyl)phthalate	21.71	149	153103	5.56	ng	# 97
87) Benzo(b)fluoranthene	24.13	252	154252	2.94	ng	# 93

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

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