

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072115\
 Data File : BG018046.D
 Acq On : 22 Jul 2015 6:43
 Operator : TP/UM
 Sample : G3027-07
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-06-1.0-1.5

Quant Time: Jul 22 07:27:58 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 22 04:14:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	53282	20.00	ng	0.00
21) Naphthalene-d8	10.33	136	255369	20.00	ng	0.00
38) Acenaphthene-d10	14.21	164	161289	20.00	ng	0.00
63) Phenanthrene-d10	16.96	188	328557	20.00	ng	0.00
75) Chrysene-d12	21.16	240	352396	20.00	ng	0.00
86) Perylene-d12	23.36	264	353178	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	431544	141.29	ng	0.00
7) Phenol-d6	6.74	99	636749	159.13	ng	0.00
23) Nitrobenzene-d5	8.70	82	384149	97.43	ng	0.00
41) 2,4,6-Tribromophenol	15.71	330	300484	183.88	ng	0.00
44) 2-Fluorobiphenyl	12.82	172	980446	105.34	ng	0.00
78) Terphenyl-d14	19.61	244	1346905	96.28	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	13.67	163	171649	15.32	ng	# 98
70) Phenanthrene	17.00	178	71155	4.20	ng	95
74) Fluoranthene	19.03	202	125231	6.75	ng	95
77) Pyrene	19.39	202	108529	5.52	ng	94
80) Benzo(a)anthracene	21.15	228	57018	3.06	ng	97
82) Chrysene	21.20	228	57567	3.25	ng	95
87) Benzo(b)fluoranthene	22.70	252	77918m	4.03	ng	
89) Benzo(a)pyrene	23.26	252	53936	3.00	ng	96

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

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