

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072115\
 Data File : BG018038.D
 Acq On : 22 Jul 2015 1:56
 Operator : TP/UM
 Sample : G3031-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-3-(2-4)

Quant Time: Jul 22 05:11:42 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	54782	20.00	ng	0.00
21) Naphthalene-d8	10.33	136	254559	20.00	ng	0.00
38) Acenaphthene-d10	14.21	164	159698	20.00	ng	0.00
63) Phenanthrene-d10	16.96	188	335754	20.00	ng	0.00
75) Chrysene-d12	21.17	240	356530	20.00	ng	0.00
86) Perylene-d12	23.36	264	356823	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	291349	92.78	ng	0.00
7) Phenol-d6	6.74	99	446394	108.50	ng	0.00
23) Nitrobenzene-d5	8.70	82	262303	66.74	ng	0.00
41) 2,4,6-Tribromophenol	15.71	330	209652	129.58	ng	0.00
44) 2-Fluorobiphenyl	12.83	172	750345	81.42	ng	0.00
78) Terphenyl-d14	19.61	244	1050223	74.20	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	13.67	163	118998	10.73	ng	99
70) Phenanthrene	17.01	178	249633	14.41	ng	98
71) Anthracene	17.09	178	67701	4.02	ng	97
74) Fluoranthene	19.03	202	346648	18.28	ng	96
77) Pyrene	19.40	202	282636	14.21	ng	96
80) Benzo(a)anthracene	21.15	228	171726	9.12	ng	98
82) Chrysene	21.20	228	149634	8.34	ng	96
83) Bis(2-ethylhexyl)phthalate	21.10	149	283118	23.09	ng	98
85) Indeno(1,2,3-cd)pyrene	25.58	276	88151	4.06	ng	# 93
87) Benzo(b)fluoranthene	22.71	252	205688m	10.54	ng	
88) Benzo(k)fluoranthene	22.74	252	64777m	3.32	ng	
89) Benzo(a)pyrene	23.27	252	137902	7.58	ng	97
91) Benzo(g,h,i)perylene	26.26	276	82159	4.52	ng	98

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

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