

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042815\
 Data File : BG016775.D
 Acq On : 28 Apr 2015 21:20
 Operator : TP/IZ
 Sample : G2022-18 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-07-0-2.0

Quant Time: Apr 29 01:47:09 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 29 00:55:31 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	71499	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	308749	20.00	ng	0.00
38) Acenaphthene-d10	14.22	164	164110	20.00	ng	0.00
63) Phenanthrene-d10	16.97	188	300803	20.00	ng	0.00
75) Chrysene-d12	21.16	240	255215	20.00	ng	0.00
86) Perylene-d12	23.36	264	251147	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	71228	16.17	ng	0.00
7) Phenol-d6	6.77	99	97930	15.84	ng	0.00
23) Nitrobenzene-d5	8.73	82	50011	10.24	ng	0.00
41) 2,4,6-Tribromophenol	15.72	330	19148	15.73	ng	0.00
44) 2-Fluorobiphenyl	12.84	172	115724	11.33	ng	0.00
78) Terphenyl-d14	19.60	244	107591	9.71	ng	0.00
Target Compounds						Qvalue
48) Acenaphthylene	13.95	152	60759	3.50	ng	97
70) Phenanthrene	17.01	178	345103	19.97	ng	100
71) Anthracene	17.10	178	56178	3.27	ng	99
72) Carbazole	17.38	167	35068	2.04	ng	97
74) Fluoranthene	19.03	202	746741	40.78	ng	98
77) Pyrene	19.40	202	727523	40.58	ng	100
80) Benzo(a)anthracene	21.14	228	357287	23.09	ng	99
82) Chrysene	21.20	228	353446	23.82	ng	98
85) Indeno(1,2,3-cd)pyrene	25.58	276	217142	13.43	ng	# 86
87) Benzo(b)fluoranthene	22.70	252	440752	28.89	ng	99
88) Benzo(k)fluoranthene	22.73	252	173060m	11.62	ng	
89) Benzo(a)pyrene	23.26	252	333608	23.21	ng	96
90) Dibenzo(a,h)anthracene	25.58	278	52261	3.74	ng	# 68
91) Benzo(a,h,i)perylene	26.26	276	198006	13.92	ng	96

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

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