

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042815\
 Data File : BG016770.D
 Acq On : 28 Apr 2015 18:26
 Operator : TP/IZ
 Sample : G1991-13DL 25X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-63SDL

Quant Time: Apr 29 01:33:12 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	68314	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	304182	20.00	ng	0.00
38) Acenaphthene-d10	14.22	164	165885	20.00	ng	0.00
63) Phenanthrene-d10	16.97	188	306146	20.00	ng	0.00
75) Chrysene-d12	21.16	240	257747	20.00	ng	0.00
86) Perylene-d12	23.36	264	248222	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	17278	4.10	ng	0.00
7) Phenol-d6	6.77	99	22579	3.82	ng	0.00
23) Nitrobenzene-d5	8.73	82	11797	2.45	ng	0.00
41) 2,4,6-Tribromophenol	15.72	330	4132	3.36	ng	0.00
44) 2-Fluorobiphenyl	12.84	172	29273	2.84	ng	0.00
78) Terphenyl-d14	19.60	244	30759	2.75	ng	0.00
Target Compounds						Qvalue
51) Acenaphthene	14.28	154	25854	2.46	ng	94
57) Fluorene	15.28	166	27915	2.15	ng	98
70) Phenanthrene	17.01	178	642198	36.51	ng	99
71) Anthracene	17.10	178	149055	8.52	ng	97
72) Carbazole	17.38	167	47843	2.73	ng	96
74) Fluoranthene	19.04	202	960697	51.55	ng	97
77) Pyrene	19.40	202	871123	48.11	ng	99
80) Benzo(a)anthracene	21.14	228	359665	23.01	ng	99
82) Chrysene	21.19	228	318560	21.26	ng	98
85) Indeno(1,2,3-cd)pyrene	25.58	276	143156	8.77	ng	# 89
87) Benzo(b)fluoranthene	22.70	252	324068m	21.49	ng	
88) Benzo(k)fluoranthene	22.73	252	127205m	8.65	ng	
89) Benzo(a)pyrene	23.26	252	258075	18.16	ng	96
90) Dibenzo(a,h)anthracene	25.58	278	38559	2.79	ng	# 64
91) Benzo(a,h,i)perylene	26.27	276	135871	9.67	ng	100

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

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