

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
 Data File : BG016751.D
 Acq On : 28 Apr 2015 4:00
 Operator : TP/IZ
 Sample : G1992-08
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TEC02

Quant Time: Apr 28 05:42:58 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 28 01:49:03 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	72933	20.00	ng	0.00
21) Naphthalene-d8	10.36	136	316617	20.00	ng	0.00
38) Acenaphthene-d10	14.22	164	177985	20.00	ng	0.00
63) Phenanthrene-d10	16.97	188	332100	20.00	ng	0.00
75) Chrysene-d12	21.16	240	301830	20.00	ng	0.00
86) Perylene-d12	23.35	264	293478	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	464349	103.32	ng	0.00
7) Phenol-d6	6.78	99	635130	100.73	ng	0.00
23) Nitrobenzene-d5	8.73	82	316045	63.07	ng	0.00
41) 2,4,6-Tribromophenol	15.72	330	135502	102.62	ng	0.00
44) 2-Fluorobiphenyl	12.84	172	692958	62.55	ng	0.00
78) Terphenyl-d14	19.60	244	668827	51.04	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	13.69	163	60347	4.53	ng	98
70) Phenanthrene	17.01	178	89318	4.68	ng	99
74) Fluoranthene	19.03	202	401210	19.85	ng	97
77) Pyrene	19.40	202	346416	16.34	ng	99
80) Benzo(a)anthracene	21.14	228	215358	11.77	ng	99
82) Chrysene	21.20	228	264104	15.05	ng	98
85) Indeno(1,2,3-cd)pyrene	25.58	276	129456	6.77	ng	# 85
87) Benzo(b)fluoranthene	22.69	252	343722	19.28	ng	97
88) Benzo(k)fluoranthene	22.74	252	114674m	6.59	ng	
89) Benzo(a)pyrene	23.26	252	210690	12.54	ng	96
91) Benzo(g,h,i)perylene	26.26	276	128965	7.76	ng	96

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

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