

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042815\
 Data File : BG016781.D
 Acq On : 29 Apr 2015 00:49
 Operator : TP/IZ
 Sample : G2021-09
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PS-18-0-2

Quant Time: Apr 29 04:56:52 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	7.57	152	72904	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	306804	20.00	ng	0.00
38) Acenaphthene-d10	14.22	164	165143	20.00	ng	0.00
63) Phenanthrene-d10	16.97	188	294508	20.00	ng	0.00
75) Chrysene-d12	21.16	240	248359	20.00	ng	0.00
86) Perylene-d12	23.35	264	245319	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	576464	128.31	ng	0.00
7) Phenol-d6	6.77	99	781741	124.03	ng	0.00
23) Nitrobenzene-d5	8.73	82	409149	84.27	ng	0.00
41) 2,4,6-Tribromophenol	15.72	330	148409	121.14	ng	0.00
44) 2-Fluorobiphenyl	12.84	172	931061	90.58	ng	0.00
78) Terphenyl-d14	19.60	244	833540	77.31	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.80	94	27807	4.19	ng	96
49) Dimethylphthalate	13.68	163	32479	2.63	ng	# 97
70) Phenanthrene	17.01	178	600065	35.47	ng	99
71) Anthracene	17.10	178	90975	5.41	ng	95
74) Fluoranthene	19.03	202	851021	47.47	ng	97
77) Pyrene	19.40	202	966514	55.39	ng	98
80) Benzo(a)anthracene	21.14	228	436255	28.97	ng	99
82) Chrysene	21.20	228	414844m	28.73	ng	
85) Indeno(1,2,3-cd)pyrene	25.57	276	160612	10.21	ng	# 86
87) Benzo(b)fluoranthene	22.70	252	438900m	29.45	ng	
88) Benzo(k)fluoranthene	22.73	252	139559m	9.60	ng	
89) Benzo(a)pyrene	23.26	252	333298	23.74	ng	98
90) Dibenzo(a,h)anthracene	25.58	278	31777m	2.33	ng	
91) Benzo(a,h,i)perylene	26.26	276	153741	11.07	ng	98

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

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