

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
 Data File : BG016735.D
 Acq On : 27 Apr 2015 14:19
 Operator : TP/IZ
 Sample : G1950-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PS-21(2-5)

Quant Time: Apr 28 01:28:19 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 00:54:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	61974	20.00	ng	0.00
21) Naphthalene-d8	10.37	136	273033	20.00	ng	0.00
38) Acenaphthene-d10	14.23	164	160908	20.00	ng	0.00
63) Phenanthrene-d10	16.98	188	312417	20.00	ng	0.00
75) Chrysene-d12	21.17	240	283429	20.00	ng	0.00
86) Perylene-d12	23.37	264	271622	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	411338	107.71	ng	0.00
7) Phenol-d6	6.78	99	573362	107.01	ng	0.00
23) Nitrobenzene-d5	8.74	82	288893	66.86	ng	0.00
41) 2,4,6-Tribromophenol	15.73	330	128790	107.89	ng	0.00
44) 2-Fluorobiphenyl	12.85	172	694858	69.38	ng	0.00
78) Terphenyl-d14	19.62	244	736370	59.85	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	13.70	163	86713	7.20	ng	99
70) Phenanthrene	17.02	178	132535	7.38	ng	98
74) Fluoranthene	19.05	202	232120	12.21	ng	95
77) Pyrene	19.41	202	221826	11.14	ng	98
80) Benzo(a)anthracene	21.16	228	149199	8.68	ng	98
82) Chrysene	21.20	228	150908	9.16	ng	97
85) Indeno(1,2,3-cd)pyrene	25.60	276	97264	5.42	ng	89
87) Benzo(b)fluoranthene	22.71	252	199750m	12.11	ng	
88) Benzo(k)fluoranthene	22.75	252	73209m	4.55	ng	
89) Benzo(a)pyrene	23.28	252	122702	7.89	ng	98
91) Benzo(g,h,i)perylene	26.28	276	93549	6.08	ng	99

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

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