

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
 Data File : BF079725.D
 Acq On : 5 Jun 2015 4:14
 Operator : TP/IZ
 Sample : G2474-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C-5

Quant Time: Jun 06 03:50:22 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 06 03:31:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	98875	20.00	ng	0.00
21) Naphthalene-d8	8.80	136	387007	20.00	ng	0.00
38) Acenaphthene-d10	10.58	164	201882	20.00	ng	0.00
63) Phenanthrene-d10	12.09	188	432065	20.00	ng	0.01
75) Chrysene-d12	15.09	240	449674	20.00	ng	0.10
86) Perylene-d12	17.05	264	447935	20.00	ng	0.14
System Monitoring Compounds						
5) 2-Fluorophenol	6.12	112	625110	108.41	ng	0.00
7) Phenol-d6	7.11	99	892990	118.96	ng	0.01
23) Nitrobenzene-d5	8.07	82	447320	70.37	ng	0.00
41) 2,4,6-Tribromophenol	11.37	330	213657	117.85	ng	0.00
44) 2-Fluorobiphenyl	9.87	172	923111	66.20	ng	0.00
78) Terphenyl-d14	13.79	244	1120739	58.72	ng	0.06
Target Compounds						Qvalue
49) Dimethylphthalate	10.26	163	124088	7.98	ng	97
70) Phenanthrene	12.11	178	293738	11.69	ng	98
71) Anthracene	12.17	178	53459	2.15	ng	98
74) Fluoranthene	13.38	202	481943	17.68	ng	96
77) Pyrene	13.65	202	423721	15.16	ng	98
80) Benzo(a)anthracene	15.06	228	213516	7.77	ng	99
82) Chrysene	15.11	228	218471	8.49	ng	97
85) Indeno(1,2,3-cd)pyrene	19.06	276	140484	4.84	ng	95
87) Benzo(b)fluoranthene	16.47	252	284911m	9.86	ng	
88) Benzo(k)fluoranthene	16.50	252	97862m	3.55	ng	
89) Benzo(a)pyrene	16.97	252	186658	7.07	ng	98
91) Benzo(g,h,i)perylene	19.67	276	134441	5.39	ng	99

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

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