

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040815\
 Data File : BF078349.D
 Acq On : 8 Apr 2015 22:28
 Operator : TP/IZ
 Sample : G1794-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OUTFALL-1

Quant Time: Apr 09 11:48:24 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 09 00:49:15 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.91	152	43363	20.00	ng	0.00
21) Naphthalene-d8	8.49	136	178062	20.00	ng	0.00
38) Acenaphthene-d10	10.65	164	88038	20.00	ng	0.00
63) Phenanthrene-d10	12.48	188	174403	20.00	ng	0.00
75) Chrysene-d12	15.75	240	178532	20.00	ng	-0.02
86) Perylene-d12	17.48	264	168840	20.00	ng	-0.09
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	179161	68.12	ng	0.02
7) Phenol-d6	6.52	99	232040	70.40	ng	0.01
23) Nitrobenzene-d5	7.61	82	128202	43.64	ng	0.00
41) 2,4,6-Tribromophenol	11.63	330	52625	72.07	ng	0.00
44) 2-Fluorobiphenyl	9.84	172	232982	37.83	ng	0.00
78) Terphenyl-d14	14.48	244	236456	29.93	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	10.35	163	16283	2.46	ng	# 98
70) Phenanthrene	12.51	178	149622	14.83	ng	99
71) Anthracene	12.57	178	23107	2.32	ng	98
72) Carbazole	12.79	167	18695	2.01	ng	98
74) Fluoranthene	13.97	202	336349	31.61	ng	97
77) Pyrene	14.25	202	264542	23.60	ng	# 91
80) Benzo(a)anthracene	15.73	228	139733	13.15	ng	96
82) Chrysene	15.78	228	144360	14.46	ng	97
85) Indeno(1,2,3-cd)pyrene	18.76	276	96733m	8.54	ng	
87) Benzo(b)fluoranthene	17.04	252	210803	20.20	ng	# 93
88) Benzo(k)fluoranthene	17.07	252	64056m	6.91	ng	
89) Benzo(a)pyrene	17.43	252	119013	13.07	ng	96
90) Dibenzo(a,h)anthracene	18.79	278	18786m	2.06	ng	
91) Benzo(g,h,i)perylene	19.13	276	86976	9.51	ng	100

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

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