

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040815\
 Data File : BF078350.D
 Acq On : 8 Apr 2015 22:56
 Operator : TP/IZ
 Sample : G1794-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OUTFALL-2

Quant Time: Apr 09 11:51:37 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	41031	20.00	ng	0.00
21) Naphthalene-d8	8.49	136	168404	20.00	ng	0.00
38) Acenaphthene-d10	10.65	164	79144	20.00	ng	0.00
63) Phenanthrene-d10	12.48	188	152090	20.00	ng	0.00
75) Chrysene-d12	15.75	240	164352	20.00	ng	-0.02
86) Perylene-d12	17.45	264	159528	20.00	ng	-0.13

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.18	112	248883	100.01	ng	0.01
7) Phenol-d6	6.52	99	324531	104.06	ng	0.01
23) Nitrobenzene-d5	7.61	82	181802	65.44	ng	0.00
41) 2,4,6-Tribromophenol	11.63	330	72424	110.34	ng	0.00
44) 2-Fluorobiphenyl	9.84	172	332333	60.02	ng	0.00
78) Terphenyl-d14	14.48	244	327124	44.98	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.53	94	10260	2.75	ng	88
31) Naphthalene	8.51	128	37659	4.30	ng	99
37) 2-Methylnaphthalene	9.38	142	37749	6.34	ng	96
49) Dimethylphthalate	10.35	163	18476	3.11	ng	99
51) Acenaphthene	10.69	154	69838	15.08	ng	98
54) Dibenzofuran	10.91	168	75704	10.70	ng	92
57) Fluorene	11.33	166	75457	13.16	ng	99
70) Phenanthrene	12.51	178	472424	53.70	ng	99
71) Anthracene	12.57	178	140896	16.25	ng	99
72) Carbazole	12.78	167	62690	7.72	ng	99
74) Fluoranthene	13.97	202	493868	53.23	ng	98
77) Pvrene	14.25	202	396013	38.38	ng	# 90
80) Benzo(a)anthracene	15.73	228	227174	23.23	ng	99
82) Chrysene	15.78	228	176820	19.25	ng	97
85) Indeno(1,2,3-cd)pyrene	18.71	276	86946m	8.34	ng	
87) Benzo(b)fluoranthene	17.01	252	206591m	20.95	ng	
88) Benzo(k)fluoranthene	17.03	252	94076m	10.74	ng	
89) Benzo(a)pyrene	17.38	252	151850	17.65	ng	# 94
90) Dibenzo(a,h)anthracene	18.73	278	19974m	2.32	ng	
91) Benzo(a,h,i)perylene	19.07	276	76884	8.90	ng	98

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

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