

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040815\
 Data File : BF078347.D
 Acq On : 8 Apr 2015 21:31
 Operator : TP/IZ
 Sample : G1793-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-2S

Quant Time: Apr 09 01:49:55 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	44877	20.00	ng	0.00
21) Naphthalene-d8	8.49	136	180204	20.00	ng	0.00
38) Acenaphthene-d10	10.65	164	86943	20.00	ng	0.00
63) Phenanthrene-d10	12.48	188	167343	20.00	ng	0.00
75) Chrysene-d12	15.77	240	176881	20.00	ng	0.00
86) Perylene-d12	17.59	264	168234	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	95932	35.25	ng	0.01
7) Phenol-d6	6.51	99	287353	84.24	ng	0.00
23) Nitrobenzene-d5	7.61	82	201469	67.77	ng	0.00
41) 2,4,6-Tribromophenol	11.63	330	15703	21.78	ng	0.00
44) 2-Fluorobiphenyl	9.84	172	359181	59.05	ng	0.00
78) Terphenyl-d14	14.48	244	387868	49.55	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.51	128	22969	2.45	ng	97
49) Dimethylphthalate	10.35	163	37231	5.70	ng	98
51) Acenaphthene	10.69	154	11126	2.19	ng	99
54) Dibenzofuran	10.91	168	21882	2.82	ng	94
57) Fluorene	11.33	166	16411	2.61	ng	# 93
70) Phenanthrene	12.51	178	236179	24.40	ng	100
71) Anthracene	12.57	178	55280	5.80	ng	96
72) Carbazole	12.79	167	25127	2.81	ng	98
74) Fluoranthene	13.97	202	214861	21.05	ng	95
77) Pyrene	14.25	202	181641	16.36	ng	# 93
80) Benzo(a)anthracene	15.76	228	100879	9.59	ng	97
82) Chrysene	15.80	228	75304	7.62	ng	97
85) Indeno(1,2,3-cd)pyrene	18.90	276	35633	3.18	ng	# 82
87) Benzo(b)fluoranthene	17.12	252	81140m	7.80	ng	
88) Benzo(k)fluoranthene	17.13	252	37482m	4.06	ng	
89) Benzo(a)pyrene	17.52	252	63803	7.03	ng	# 88
91) Benzo(a,h,i)perylene	19.27	276	36151	3.97	ng	96

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

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