

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040915\
 Data File : BF078393.D
 Acq On : 10 Apr 2015 9:42
 Operator : TP/IZ
 Sample : G1793-09DL 10X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-6DADL

Manual Integrations
 APPROVED

apatel
 4/10/2015 5:13:50 PM

Quant Time: Apr 10 13:02:38 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 11:29:31 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	35747	20.00	ng	0.00
21) Naphthalene-d8	8.46	136	142985	20.00	ng	0.00
38) Acenaphthene-d10	10.62	164	73163	20.00	ng	-0.01
63) Phenanthrene-d10	12.45	188	139473	20.00	ng	-0.01
75) Chrysene-d12	15.73	240	139082	20.00	ng	-0.01
86) Perylene-d12	17.49	264	145172	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	8305	3.83	ng	-0.01
7) Phenol-d6	6.49	99	13510	4.97	ng	0.01
23) Nitrobenzene-d5	7.58	82	7684	3.26	ng	0.00
41) 2,4,6-Tribromophenol	11.61	330	4244	6.99	ng	-0.01
44) 2-Fluorobiphenyl	9.81	172	23243	4.54	ng	0.00
78) Terphenyl-d14	14.44	244	30346	4.93	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	8.49	128	704760	94.70	ng	99
37) 2-Methylnaphthalene	9.34	142	178453	35.32	ng	98
45) 1,1'-Biphenyl	9.93	154	66307	11.08	ng	99
48) Acenaphthylene	10.45	152	42084	5.53	ng	# 94
51) Acenaphthene	10.66	154	151323	35.35	ng	98
54) Dibenzofuran	10.88	168	99580	15.23	ng	98
57) Fluorene	11.30	166	167135	31.54	ng	100
70) Phenanthrene	12.48	178	681771	84.51	ng	99
71) Anthracene	12.55	178	244993	30.82	ng	100
72) Carbazole	12.76	167	25935	3.48	ng	99
74) Fluoranthene	13.95	202	384104	45.14	ng	90
77) Pyrene	14.23	202	375925	43.06	ng	99
80) Benzo(a)anthracene	15.72	228	172824	20.88	ng	97
82) Chrysene	15.76	228	131639	16.93	ng	99
85) Indeno(1,2,3-cd)pyrene	18.76	276	47410	5.37	ng	# 88
87) Benzo(b)fluoranthene	17.04	252	135677m	15.12	ng	
88) Benzo(k)fluoranthene	17.07	252	37687m	4.73	ng	
89) Benzo(a)pyrene	17.43	252	121490	15.52	ng	# 95
91) Benzo(a,h,i)perylene	19.13	276	44537m	5.67	ng	

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

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