

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040915\
 Data File : BF078396.D
 Acq On : 10 Apr 2015 11:08
 Operator : TP/IZ
 Sample : G1794-02DL 20X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DBOXDL

Manual Integrations
 APPROVED

apatel
 4/10/2015 5:14:02 PM

Quant Time: Apr 10 13:16:36 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	41350	20.00	ng	0.00
21) Naphthalene-d8	8.46	136	179187	20.00	ng	0.00
38) Acenaphthene-d10	10.62	164	86999	20.00	ng	-0.01
63) Phenanthrene-d10	12.45	188	162919	20.00	ng	-0.01
75) Chrysene-d12	15.72	240	166279	20.00	ng	-0.02
86) Perylene-d12	17.47	264	170153	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	8020	3.20	ng	-0.01
7) Phenol-d6	6.50	99	12127	3.86	ng	0.02
23) Nitrobenzene-d5	7.58	82	6453	2.18	ng	0.00
41) 2,4,6-Tribromophenol	11.60	330	2599	3.60	ng	-0.02
44) 2-Fluorobiphenyl	9.81	172	16801	2.76	ng	0.00
78) Terphenyl-d14	14.44	244	18051	2.45	ng	-0.02
Target Compounds						Qvalue
70) Phenanthrene	12.48	178	183035	19.42	ng	100
71) Anthracene	12.54	178	29928	3.22	ng	97
72) Carbazole	12.76	167	23593	2.71	ng	100
74) Fluoranthene	13.95	202	470517	47.34	ng	90
77) Pyrene	14.23	202	389576	37.32	ng	98
80) Benzo(a)anthracene	15.71	228	203692	20.59	ng	98
82) Chrysene	15.76	228	241829	26.02	ng	98
85) Indeno(1,2,3-cd)pyrene	18.75	276	167328m	15.86	ng	
87) Benzo(b)fluoranthene	17.03	252	327887m	31.18	ng	
88) Benzo(k)fluoranthene	17.04	252	131805m	14.10	ng	
89) Benzo(a)pyrene	17.40	252	206933	22.55	ng	# 91
90) Dibenzo(a,h)anthracene	18.77	278	33820m	3.68	ng	
91) Benzo(a,h,i)perylene	19.11	276	160067m	17.37	ng	

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

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