

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040915\
 Data File : BF078397.D
 Acq On : 10 Apr 2015 11:37
 Operator : TP/IZ
 Sample : G1794-04RE
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OUTFALL-1RE

Manual Integrations
 APPROVED

apatel
 4/10/2015 5:14:07 PM

Quant Time: Apr 10 13:21:25 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	47040	20.00	ng	0.00
21) Naphthalene-d8	8.46	136	198139	20.00	ng	0.00
38) Acenaphthene-d10	10.63	164	98778	20.00	ng	-0.01
63) Phenanthrene-d10	12.45	188	184140	20.00	ng	-0.01
75) Chrysene-d12	15.71	240	188478	20.00	ng	-0.03
86) Perylene-d12	17.43	264	185192	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.15	112	200645	70.33	ng	-0.02
7) Phenol-d6	6.49	99	258033	72.17	ng	0.01
23) Nitrobenzene-d5	7.58	82	144446	44.19	ng	0.00
41) 2,4,6-Tribromophenol	11.60	330	53783	65.65	ng	-0.02
44) 2-Fluorobiphenyl	9.81	172	262108	37.93	ng	0.00
78) Terphenyl-d14	14.44	244	248676	29.81	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	10.33	163	17037	2.30	ng	99
70) Phenanthrene	12.48	178	159621	14.99	ng	99
71) Anthracene	12.55	178	24279	2.31	ng	98
74) Fluoranthene	13.94	202	359096	31.97	ng	98
77) Pyrene	14.21	202	284302	24.03	ng	# 92
80) Benzo(a)anthracene	15.70	228	153275	13.67	ng	96
82) Chrysene	15.75	228	149123	14.15	ng	98
85) Indeno(1,2,3-cd)pyrene	18.68	276	96574m	8.08	ng	
87) Benzo(b)fluoranthene	16.99	252	251846m	22.00	ng	
88) Benzo(k)fluoranthene	17.03	252	52029m	5.12	ng	
89) Benzo(a)pyrene	17.37	252	133976	13.41	ng	# 96
90) Dibenzo(a,h)anthracene	18.71	278	20287m	2.03	ng	
91) Benzo(a,h,i)perylene	19.05	276	95060m	9.48	ng	

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

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