

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040915\
 Data File : BF078390.D
 Acq On : 10 Apr 2015 8:17
 Operator : TP/IZ
 Sample : G1793-07DL 5X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-5DDL

Manual Integrations
 APPROVED

apatel
 4/10/2015 5:13:36 PM

Quant Time: Apr 10 12:38:51 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	39473	20.00	ng	0.00
21) Naphthalene-d8	8.48	136	160404	20.00	ng	0.01
38) Acenaphthene-d10	10.63	164	83614	20.00	ng	-0.01
63) Phenanthrene-d10	12.45	188	150854	20.00	ng	-0.01
75) Chrysene-d12	15.72	240	146739	20.00	ng	-0.02
86) Perylene-d12	17.45	264	149496	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	39869	16.65	ng	-0.01
7) Phenol-d6	6.49	99	51715	17.24	ng	0.01
23) Nitrobenzene-d5	7.58	82	25285	9.55	ng	0.00
41) 2,4,6-Tribromophenol	11.61	330	10111	14.58	ng	-0.01
44) 2-Fluorobiphenyl	9.81	172	60363	10.32	ng	0.00
78) Terphenyl-d14	14.44	244	62579	9.64	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	8.51	128	3358895	402.32	ng	95
37) 2-Methylnaphthalene	9.36	142	812992	143.43	ng	96
45) 1,1'-Biphenyl	9.94	154	356831	52.17	ng	99
48) Acenaphthylene	10.45	152	311743	35.88	ng	# 97
51) Acenaphthene	10.67	154	516606	105.59	ng	99
54) Dibenzofuran	10.89	168	125191	16.75	ng	93
57) Fluorene	11.31	166	546248	90.18	ng	98
70) Phenanthrene	12.50	178	2024120	231.96	ng	97
71) Anthracene	12.56	178	743679	86.49	ng	100
72) Carbazole	12.76	167	36797	4.57	ng	96
74) Fluoranthene	13.95	202	916074	99.55	ng	98
77) Pyrene	14.24	202	1263086	137.12	ng	97
80) Benzo(a)anthracene	15.71	228	438715m	50.25	ng	
82) Chrysene	15.76	228	319845	38.99	ng	98
85) Indeno(1,2,3-cd)pyrene	18.71	276	117676	12.64	ng	# 88
87) Benzo(b)fluoranthene	17.01	252	288643m	31.24	ng	
88) Benzo(k)fluoranthene	17.04	252	79921m	9.73	ng	
89) Benzo(a)pyrene	17.38	252	326071	40.44	ng	# 88
90) Dibenzo(a,h)anthracene	18.73	278	27955m	3.46	ng	
91) Benzo(a,h,i)perylene	19.07	276	141675m	17.50	ng	

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

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