

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042115\
 Data File : BF078694.D
 Acq On : 21 Apr 2015 20:36
 Operator : TP/IZ
 Sample : G1938-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-65S

Manual Integrations
 APPROVED

apatel
 4/22/2015 5:22:29 PM

Quant Time: Apr 22 03:35:51 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 22 00:52:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.82	152	32520	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	134571	20.00	ng	-0.01
38) Acenaphthene-d10	11.17	164	68920	20.00	ng	-0.01
63) Phenanthrene-d10	13.90	188	134136	20.00	ng	-0.01
75) Chrysene-d12	17.77	240	145715	20.00	ng	-0.01
86) Perylene-d12	19.50	264	156670	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	3.84	112	176144	89.31	ng	0.01
7) Phenol-d6	5.32	99	230630	93.30	ng	0.00
23) Nitrobenzene-d5	6.75	82	133182	59.99	ng	-0.01
41) 2,4,6-Tribromophenol	12.64	330	57729	101.00	ng	-0.01
44) 2-Fluorobiphenyl	9.99	172	282179	58.52	ng	-0.01
78) Terphenyl-d14	16.43	244	327876	50.84	ng	0.00
Target Compounds						Qvalue
48) Acenaphthylene	10.90	152	27670	3.86	ng	97
49) Dimethylphthalate	10.80	163	11941	2.31	ng	# 95
51) Acenaphthene	11.22	154	15565	3.86	ng	98
54) Dibenzofuran	11.55	168	14652	2.38	ng	99
57) Fluorene	12.18	166	23426	4.69	ng	98
70) Phenanthrene	13.94	178	362241	46.69	ng	100
71) Anthracene	14.03	178	92686	12.12	ng	99
72) Carbazole	14.37	167	27174	3.79	ng	95
74) Fluoranthene	15.84	202	645373	78.87	ng	94
77) Pyrene	16.15	202	566827	61.97	ng	# 92
80) Benzo(a)anthracene	17.76	228	360787	41.61	ng	98
82) Chrysene	17.81	228	272553	33.46	ng	97
85) Indeno(1,2,3-cd)pyrene	20.69	276	196032m	21.21	ng	
87) Benzo(b)fluoranthene	19.07	252	416013m	42.96	ng	
88) Benzo(k)fluoranthene	19.10	252	133025m	15.46	ng	
89) Benzo(a)pyrene	19.44	252	303221	35.88	ng	96
90) Dibenzo(a,h)anthracene	20.71	278	46926m	5.55	ng	
91) Benzo(g,h,i)perylene	21.01	276	179248	21.13	ng	100

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

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