

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF021918\
 Data File : BF103098.D
 Acq On : 19 Feb 2018 23:09
 Operator : SJ/JU
 Sample : J1600-08
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 A2 SB44 BOT-S 8

Manual Integrations
 APPROVED

Sohil
 2/20/2018 10:20:53 AM

Quant Time: Feb 20 01:20:10 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 16 14:34:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	177119	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	741460	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	314747	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	461239	20.00	ng	0.00
75) Chrysene-d12	14.06	240	237705	20.00	ng	0.00
86) Perylene-d12	15.55	264	285488	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1124853	96.45	ng	0.00
7) Phenol-d6	6.51	99	1348448	96.02	ng	0.00
23) Nitrobenzene-d5	7.44	82	876491	82.00	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	240047	94.75	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1201706	63.35	ng	0.00
78) Terphenyl-d14	12.99	244	712745	71.00	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.52	94	31045	2.063	ng	95
31) Naphthalene	8.18	128	297364	8.209	ng	100
37) 2-Methylnaphthalene	8.87	142	69860	2.945	ng	97
48) Acenaphthylene	9.77	152	549305	16.946	ng	99
49) Dimethylphthalate	9.62	163	127517	5.196	ng	99
54) Dibenzofuran	10.12	168	172926	6.330	ng	97
57) Fluorene	10.46	166	384945	19.366	ng	99
70) Phenanthrene	11.44	178	1939735	75.511	ng	99
71) Anthracene	11.49	178	840358	32.164	ng	99
72) Carbazole	11.63	167	132068	5.160	ng	99
74) Fluoranthene	12.64	202	2812391	108.817	ng	99
77) Pyrene	12.87	202	2192027	117.599	ng	99
80) Benzo(a)anthracene	14.05	228	1460707m	97.710	ng	
82) Chrysene	14.09	228	1316607m	87.367	ng	
85) Indeno(1,2,3-cd)pyrene	17.07	276	664272	53.148	ng	95
87) Benzo(b)fluoranthene	15.12	252	1770933m	95.678	ng	
88) Benzo(k)fluoranthene	15.14	252	549534m	31.073	ng	
89) Benzo(a)pyrene	15.50	252	1207871	72.499	ng	95
90) Dibenzo(a,h)anthracene	17.07	278	198666m	14.691	ng	
91) Benzo(g,h,i)perylene	17.53	276	586657	44.323	ng	96

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

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