

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030918\
 Data File : BF103557.D
 Acq On : 9 Mar 2018 15:19
 Operator : SJ/JU
 Sample : J1853-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP1-WC

Manual Integrations
 APPROVED

Sohil
 3/12/2018 3:50:13 PM

Quant Time: Mar 09 17:30:44 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 09 11:47:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	157132	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	649730	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	263934	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	367165	20.00	ng	0.00
75) Chrysene-d12	14.06	240	229598	20.00	ng	0.00
86) Perylene-d12	15.58	264	187040	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	1028213	98.97	ng	0.00
7) Phenol-d6	6.50	99	1297152	102.03	ng	0.00
23) Nitrobenzene-d5	7.43	82	821065	72.17	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	164596	73.68	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	1011909	61.80	ng	0.00
78) Terphenyl-d14	12.99	244	629288	65.89	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.17	128	83082	2.612	ng	98
37) 2-Methylnaphthalene	8.86	142	101132	4.919	ng	97
51) Acenaphthene	9.94	154	51057	3.252	ng	95
54) Dibenzofuran	10.12	168	51720	2.400	ng	# 92
57) Fluorene	10.46	166	61310	3.340	ng	# 98
70) Phenanthrene	11.43	178	466829	23.103	ng	99
71) Anthracene	11.48	178	154676	7.465	ng	97
72) Carbazole	11.65	167	47706	2.312	ng	# 95
73) Di-n-butylphthalate	11.95	149	804248	31.149	ng	99
74) Fluoranthene	12.62	202	494630	24.360	ng	99
77) Pyrene	12.86	202	449751	25.125	ng	98
80) Benzo(a)anthracene	14.05	228	226403	15.198	ng	97
82) Chrysene	14.09	228	190856	13.195	ng	95
85) Indeno(1,2,3-cd)pyrene	17.14	276	80058	6.991	ng	# 91
87) Benzo(b)fluoranthene	15.13	252	199122m	16.192	ng	
88) Benzo(k)fluoranthene	15.16	252	78961m	6.819	ng	
89) Benzo(a)pyrene	15.52	252	156223	14.226	ng	# 87
91) Benzo(g,h,i)perylene	17.62	276	94067	11.342	ng	# 90

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

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