

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF033018\
 Data File : BF104117.D
 Acq On : 31 Mar 2018 2:10
 Operator : JU/SJ
 Sample : J2142-07
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GB-RT-4252

Manual Integrations
 APPROVED

Sohil
 4/2/2018 5:06:11 PM

Quant Time: Mar 31 04:30:11 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 30 17:20:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	109568	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	447733	20.00	ng	0.00
38) Acenaphthene-d10	9.99	164	191246	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	298141	20.00	ng	0.00
75) Chrysene-d12	14.14	240	235157	20.00	ng	0.00
86) Perylene-d12	15.69	264	201064	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	649571	94.54	ng	0.01
7) Phenol-d6	6.59	99	874107	103.41	ng	0.00
23) Nitrobenzene-d5	7.52	82	505104	68.99	ng	0.00
41) 2,4,6-Tribromophenol	10.79	330	173589	81.52	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	791522	65.26	ng	0.00
78) Terphenyl-d14	13.07	244	634655	62.32	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.70	163	65991	4.370	ng	100
57) Fluorene	10.54	166	28799	2.107	ng	# 98
70) Phenanthrene	11.52	178	393615	24.385	ng	99
71) Anthracene	11.57	178	138871	8.236	ng	97
74) Fluoranthene	12.71	202	747414	43.561	ng	98
77) Pyrene	12.94	202	618383	36.581	ng	99
80) Benzo(a)anthracene	14.13	228	340145	23.117	ng	98
82) Chrysene	14.17	228	279051	19.363	ng	97
85) Indeno(1,2,3-cd)pyrene	17.26	276	118748	7.951	ng	# 87
87) Benzo(b)fluoranthene	15.23	252	296714m	24.248	ng	
88) Benzo(k)fluoranthene	15.25	252	140656m	12.202	ng	
89) Benzo(a)pyrene	15.62	252	223905	19.745	ng	97
90) Dibenzo(a,h)anthracene	17.27	278	24220	2.310	ng	96
91) Benzo(a,h,i)perylene	17.74	276	100974	9.616	ng	96

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

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