

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF022818\
 Data File : BF103315.D
 Acq On : 1 Mar 2018 1:11
 Operator : SJ/JU
 Sample : J1720-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-31

Manual Integrations
 APPROVED

Sohil

3/1/2018 11:43:41 AM

Quant Time: Mar 01 03:03:55 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 17:05:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	144167	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	606494	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	247818	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	351955	20.00	ng	0.00
75) Chrysene-d12	14.06	240	254469	20.00	ng	0.00
86) Perylene-d12	15.56	264	201519	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1029221	107.97	ng	0.00
7) Phenol-d6	6.51	99	1245217	106.76	ng	0.00
23) Nitrobenzene-d5	7.44	82	803826	75.69	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	172044	82.02	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1091555	73.44	ng	0.00
78) Terphenyl-d14	12.99	244	704880	66.59	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	94671	4.762	ng	Qvalue 100
70) Phenanthrene	11.43	178	201630	10.410	ng	99
71) Anthracene	11.48	178	44325	2.232	ng	98
74) Fluoranthene	12.62	202	317469	16.311	ng	99
77) Pyrene	12.86	202	302671	15.256	ng	99
80) Benzo(a)anthracene	14.04	228	146486m	8.872	ng	
82) Chrysene	14.08	228	150621m	9.395	ng	
85) Indeno(1,2,3-cd)pyrene	17.09	276	57374	4.521	ng	96
87) Benzo(b)fluoranthene	15.12	252	159533m	12.041	ng	
88) Benzo(k)fluoranthene	15.14	252	61518m	4.931	ng	
89) Benzo(a)pyrene	15.49	252	107581	9.092	ng	# 93
91) Benzo(a,h,i)perylene	17.55	276	56867	6.364	ng	93

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

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