

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
 Data File : BF103480.D
 Acq On : 7 Mar 2018 2:20
 Operator : SJ/JU
 Sample : J1810-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MHA-WCD

Manual Integrations
 APPROVED

Sohil
 3/7/2018 12:17:47 PM

Quant Time: Mar 07 06:56:00 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 06 11:52:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	114191	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	473490	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	183040	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	279140	20.00	ng	0.00
75) Chrysene-d12	14.06	240	200250	20.00	ng	0.00
86) Perylene-d12	15.57	264	155930	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	884669	117.17	ng	0.00
7) Phenol-d6	6.51	99	1054736	114.16	ng	0.00
23) Nitrobenzene-d5	7.44	82	679868	82.00	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	123566	79.75	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	779575	70.40	ng	0.00
78) Terphenyl-d14	12.99	244	527195	63.29	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	91873	6.257	ng	100
70) Phenanthrene	11.43	178	34898	2.272	ng	97
74) Fluoranthene	12.63	202	71636	4.641	ng	98
77) Pyrene	12.86	202	66766	4.276	ng	99
87) Benzo(b)fluoranthene	15.13	252	20817m	2.030	ng	

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

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