

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
 Data File : BF103451.D
 Acq On : 6 Mar 2018 13:10
 Operator : SJ/JU
 Sample : J1723-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-23

Manual Integrations
 APPROVED

Sohil
 3/7/2018 12:14:47 PM

Quant Time: Mar 06 14:56:44 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	125936	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	527180	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	223208	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	352437	20.00	ng	0.00
75) Chrysene-d12	14.06	240	210024	20.00	ng	0.00
86) Perylene-d12	15.58	264	188546	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	635468	76.32	ng	0.00
7) Phenol-d6	6.51	99	778969	76.45	ng	0.00
23) Nitrobenzene-d5	7.44	82	483289	52.35	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	110750	58.62	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	716231	49.84	ng	0.00
78) Terphenyl-d14	12.99	244	480008	54.94	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	43058	2.405	ng	99
70) Phenanthrene	11.43	178	130299	6.718	ng	98
74) Fluoranthene	12.63	202	215632	11.063	ng	98
77) Pyrene	12.86	202	213769	13.055	ng	99
80) Benzo(a)anthracene	14.05	228	96701	7.096	ng	97
82) Chrysene	14.09	228	89942	6.798	ng	97
85) Indeno(1,2,3-cd)pyrene	17.13	276	39831	3.802	ng	96
87) Benzo(b)fluoranthene	15.13	252	110274m	8.895	ng	
88) Benzo(k)fluoranthene	15.16	252	42963m	3.680	ng	
89) Benzo(a)pyrene	15.52	252	75502	6.820	ng	# 91
91) Benzo(a,h,i)perylene	17.60	276	36445	4.359	ng	# 93

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

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