

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030118\
 Data File : BF103335.D
 Acq On : 1 Mar 2018 14:35
 Operator : SJ/JU
 Sample : J1721-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-41

Manual Integrations
 APPROVED

Sohil
 3/2/2018 3:30:34 PM

Quant Time: Mar 02 09:41:01 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	150697	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	560782	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	199399	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	305096	20.00	ng	0.00
75) Chrysene-d12	14.06	240	250944	20.00	ng	0.00
86) Perylene-d12	15.57	264	168980	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	855879	85.90	ng	0.01
7) Phenol-d6	6.51	99	992179	81.38	ng	0.00
23) Nitrobenzene-d5	7.44	82	653710	66.57	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	115288	68.31	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	821391	67.53	ng	0.00
78) Terphenyl-d14	12.99	244	592277	56.74	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.63	163	77307	4.833	ng	98
70) Phenanthrene	11.43	178	141442	8.424	ng	98
71) Anthracene	11.49	178	49740	2.889	ng	99
74) Fluoranthene	12.63	202	311433	18.458	ng	98
77) Pyrene	12.86	202	313006	15.998	ng	99
80) Benzo(a)anthracene	14.05	228	159516	9.797	ng	96
82) Chrysene	14.09	228	138890	8.785	ng	98
85) Indeno(1,2,3-cd)pyrene	17.10	276	48131	3.846	ng	96
87) Benzo(b)fluoranthene	15.13	252	141617m	12.746	ng	
88) Benzo(k)fluoranthene	15.16	252	49272m	4.710	ng	
89) Benzo(a)pyrene	15.51	252	92927	9.366	ng	# 86
91) Benzo(a,h,i)perylene	17.57	276	50299	6.713	ng	92

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

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