

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
 Data File : BF103458.D
 Acq On : 6 Mar 2018 16:17
 Operator : SJ/JU
 Sample : J1723-11
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-4

Manual Integrations
 APPROVED

Sohil
 3/7/2018 12:15:19 PM

Quant Time: Mar 06 17:23:49 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	115799	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	497972	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	197563	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	288216	20.00	ng	0.00
75) Chrysene-d12	14.10	240	202355	20.00	ng	0.03
86) Perylene-d12	15.66	264	148358	20.00	ng	0.09
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	750795	98.06	ng	0.00
7) Phenol-d6	6.51	99	948090	101.20	ng	0.00
23) Nitrobenzene-d5	7.44	82	559787	64.20	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	125719	75.18	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	751324	61.19	ng	0.00
78) Terphenyl-d14	12.99	244	504059	59.88	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.62	163	49964	3.153	ng	98
70) Phenanthrene	11.43	178	164415	10.366	ng	100
71) Anthracene	11.49	178	52258	3.213	ng	97
74) Fluoranthene	12.63	202	308290	19.342	ng	98
77) Pyrene	12.86	202	290842	18.435	ng	99
80) Benzo(a)anthracene	14.09	228	150060	11.429	ng	98
82) Chrysene	14.12	228	134379	10.541	ng	98
85) Indeno(1,2,3-cd)pyrene	17.22	276	47665	4.723	ng	98
87) Benzo(b)fluoranthene	15.22	252	140727m	14.427	ng	
88) Benzo(k)fluoranthene	15.24	252	48303m	5.259	ng	
89) Benzo(a)pyrene	15.60	252	92351	10.602	ng	# 89
91) Benzo(a,h,i)perylene	17.70	276	48796	7.418	ng	# 88

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

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