

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
 Data File : BF103423.D
 Acq On : 5 Mar 2018 21:49
 Operator : SJ/JU
 Sample : J1723-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-3

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:23:28 PM

Quant Time: Mar 06 00:58:44 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 05 13:40:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	130625	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	538446	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	216143	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	339795	20.00	ng	0.00
75) Chrysene-d12	14.07	240	206572	20.00	ng	0.00
86) Perylene-d12	15.59	264	158346	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	969547	112.26	ng	0.00
7) Phenol-d6	6.52	99	1152117	109.02	ng	0.00
23) Nitrobenzene-d5	7.44	82	694564	73.67	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	154637	84.52	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	885699	67.09	ng	0.00
78) Terphenyl-d14	13.00	244	616530	71.75	ng	0.00
Target Compounds						
10) Phenol	6.53	94	95072	7.749	ng	86
49) Dimethylphthalate	9.62	163	101210	5.837	ng	# 99
70) Phenanthrene	11.43	178	161131	8.617	ng	99
71) Anthracene	11.49	178	56683	2.956	ng	99
74) Fluoranthene	12.63	202	297655	15.840	ng	97
77) Pyrene	12.86	202	277634	17.238	ng	99
80) Benzo(a)anthracene	14.06	228	136230	10.164	ng	98
82) Chrysene	14.09	228	122805	9.436	ng	98
85) Indeno(1,2,3-cd)pyrene	17.14	276	51685	5.017	ng	93
87) Benzo(b)fluoranthene	15.14	252	142673m	13.704	ng	
88) Benzo(k)fluoranthene	15.16	252	45641m	4.655	ng	
89) Benzo(a)pyrene	15.52	252	93295	10.035	ng	# 88
91) Benzo(a,h,i)perylene	17.62	276	50552	7.200	ng	96

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

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