

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030818\
 Data File : BF103528.D
 Acq On : 8 Mar 2018 20:29
 Operator : SJ/JU
 Sample : J1823-01 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MHB-WCA

Manual Integrations
 APPROVED

Sohil
 3/9/2018 11:39:42 AM

Quant Time: Mar 09 00:57:03 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 08 14:20:19 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	122565	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	516036	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	195211	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	294555	20.00	ng	0.00
75) Chrysene-d12	14.06	240	217899	20.00	ng	0.00
86) Perylene-d12	15.57	264	158137	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	124238	15.33	ng	0.01
7) Phenol-d6	6.52	99	187579	18.92	ng	0.01
23) Nitrobenzene-d5	7.45	82	114869	12.71	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	20293	12.28	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	199523	12.74	ng	0.00
78) Terphenyl-d14	12.98	244	139445	15.38	ng	0.00
Target Compounds						
70) Phenanthrene	11.43	178	123479	7.617	ng	97
71) Anthracene	11.49	178	41885	2.520	ng	97
73) Di-n-butylphthalate	11.95	149	58056	2.803	ng	100
74) Fluoranthene	12.62	202	172293	10.577	ng	98
77) Pyrene	12.86	202	146944	8.649	ng	99
80) Benzo(a)anthracene	14.05	228	54165m	3.831	ng	
82) Chrysene	14.09	228	55720	4.059	ng	98
87) Benzo(b)fluoranthene	15.13	252	47502m	4.569	ng	
89) Benzo(a)pyrene	15.50	252	31678	3.412	ng	# 66
91) Benzo(g,h,i)perylene	17.60	276	16012	2.284	ng	# 76

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

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