

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030118\
 Data File : BF103328.D
 Acq On : 1 Mar 2018 11:29
 Operator : SJ/JU
 Sample : J1710-04 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-201

Manual Integrations
 APPROVED

Sohil
 3/2/2018 3:30:18 PM

Quant Time: Mar 01 18:17:19 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	146271	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	618017	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	271659	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	486197	20.00	ng	0.00
75) Chrysene-d12	14.06	240	257924	20.00	ng	0.00
86) Perylene-d12	15.56	264	247785	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	264067	27.30	ng	0.00
7) Phenol-d6	6.50	99	328723	27.78	ng	0.00
23) Nitrobenzene-d5	7.43	82	184992	17.09	ng	-0.01
41) 2,4,6-Tribromophenol	10.71	330	54574	23.73	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	303161	14.21	ng	0.00
78) Terphenyl-d14	12.99	244	236779	22.07	ng	0.00
Target Compounds						
70) Phenanthrene	11.43	178	389986	14.575	ng	99
71) Anthracene	11.48	178	97788	3.564	ng	100
74) Fluoranthene	12.62	202	376215	13.992	ng	98
77) Pyrene	12.85	202	363647	18.083	ng	99
80) Benzo(a)anthracene	14.04	228	117818	7.040	ng	99
82) Chrysene	14.08	228	104219	6.414	ng	98
85) Indeno(1,2,3-cd)pyrene	17.08	276	50873	3.955	ng	94
87) Benzo(b)fluoranthene	15.12	252	105304m	6.464	ng	
88) Benzo(k)fluoranthene	15.14	252	40194m	2.620	ng	
89) Benzo(a)pyrene	15.49	252	84854	5.833	ng	# 92
91) Benzo(g,h,i)perylene	17.55	276	56517	5.144	ng	95

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

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