

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030118\
 Data File : BF103327.D
 Acq On : 1 Mar 2018 11:02
 Operator : SJ/JU
 Sample : J1710-07 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-215

Manual Integrations
 APPROVED

Sohil
 3/2/2018 3:30:16 PM

Quant Time: Mar 01 12:58:55 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	149625	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	643566	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	275264	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	447236	20.00	ng	0.00
75) Chrysene-d12	14.06	240	238815	20.00	ng	0.00
86) Perylene-d12	15.55	264	246607	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	212171	21.45	ng	0.00
7) Phenol-d6	6.50	99	263293	21.75	ng	0.00
23) Nitrobenzene-d5	7.43	82	170915	15.17	ng	-0.01
41) 2,4,6-Tribromophenol	10.71	330	42848	18.39	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	254917	11.24	ng	0.00
78) Terphenyl-d14	12.99	244	175292	17.64	ng	0.00
Target Compounds						Qvalue
70) Phenanthrene	11.43	178	312777	12.708	ng	99
71) Anthracene	11.48	178	82153	3.255	ng	99
74) Fluoranthene	12.62	202	328622	13.287	ng	99
77) Pyrene	12.85	202	274285	14.731	ng	99
80) Benzo(a)anthracene	14.04	228	92364	5.961	ng	98
82) Chrysene	14.08	228	86804	5.769	ng	97
85) Indeno(1,2,3-cd)pyrene	17.08	276	50920	4.275	ng	96
87) Benzo(b)fluoranthene	15.11	252	100238m	6.182	ng	
88) Benzo(k)fluoranthene	15.14	252	33333m	2.183	ng	
89) Benzo(a)pyrene	15.49	252	76949	5.314	ng	# 95
91) Benzo(g,h,i)perylene	17.54	276	54602	4.993	ng	96

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

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