

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF021618\
 Data File : BF103058.D
 Acq On : 16 Feb 2018 17:53
 Operator : SJ/JU
 Sample : J1401-03 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SA-06-021518-B

Manual Integrations
 APPROVED

Sohil
 2/19/2018 11:09:43 AM

Quant Time: Feb 16 23:32:38 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 16 14:34:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	169057	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	659490	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	262789	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	398757	20.00	ng	0.00
75) Chrysene-d12	14.04	240	297433	20.00	ng	-0.01
86) Perylene-d12	15.54	264	240623	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	230399	20.70	ng	0.00
7) Phenol-d6	6.50	99	284290	21.21	ng	-0.01
23) Nitrobenzene-d5	7.43	82	160040	16.83	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	40111	18.96	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	281047	17.75	ng	-0.01
78) Terphenyl-d14	12.98	244	203645	16.21	ng	-0.01
Target Compounds						Qvalue
70) Phenanthrene	11.43	178	94722	4.265	ng	99
74) Fluoranthene	12.62	202	148519	6.647	ng	98
77) Pyrene	12.84	202	157168	6.739	ng	99
80) Benzo(a)anthracene	14.03	228	60140	3.215	ng	95
82) Chrysene	14.07	228	59697	3.166	ng	97
87) Benzo(b)fluoranthene	15.10	252	54330m	3.483	ng	
89) Benzo(a)pyrene	15.47	252	42695	3.040	ng	# 84
91) Benzo(g,h,i)perylene	17.50	276	24543	2.200	ng	# 91

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

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