

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF040218\
 Data File : BF104134.D
 Acq On : 2 Apr 2018 14:48
 Operator : JU/SJ
 Sample : J1757-09 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-03-033018-B

Manual Integrations
 APPROVED

Sohil

4/3/2018 2:27:03 PM

Quant Time: Apr 03 09:30:15 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 02 13:40:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	99747	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	433077	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	203350	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	331111	20.00	ng	0.00
75) Chrysene-d12	14.14	240	233886	20.00	ng	0.00
86) Perylene-d12	15.69	264	258383	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	106693m	17.06	ng	0.00
7) Phenol-d6	6.60	99	152825m	19.86	ng	0.00
23) Nitrobenzene-d5	7.53	82	93268m	13.17	ng	0.00
41) 2,4,6-Tribromophenol	10.79	330	35769	15.80	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	185813	14.41	ng	0.00
78) Terphenyl-d14	13.07	244	140390	13.86	ng	0.00
Target Compounds						
70) Phenanthrene	11.52	178	141523	7.895	ng	98
71) Anthracene	11.57	178	47960m	2.561	ng	
74) Fluoranthene	12.72	202	102073	5.357	ng	96
77) Pyrene	12.94	202	89877	5.346	ng	99
80) Benzo(a)anthracene	14.13	228	37394	2.555	ng	95
82) Chrysene	14.17	228	35253m	2.459	ng	
87) Benzo(b)fluoranthene	15.23	252	31788m	2.022	ng	
89) Benzo(a)pyrene	15.62	252	29872	2.050	ng	# 85

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

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