

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF040218\
 Data File : BF104136.D
 Acq On : 2 Apr 2018 15:41
 Operator : JU/SJ
 Sample : J2099-01 10X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-102(1-2)

Manual Integrations
 APPROVED

Sohil
 4/3/2018 2:27:09 PM

Quant Time: Apr 02 17:30:29 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	112025	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	484324	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	223581	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	344978	20.00	ng	0.00
75) Chrysene-d12	14.17	240	221516	20.00	ng	0.02
86) Perylene-d12	15.74	264	257382	20.00	ng	0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	61863m	8.81	ng	0.02
7) Phenol-d6	6.60	99	76849m	8.89	ng	0.00
23) Nitrobenzene-d5	7.53	82	46305m	5.85	ng	0.00
41) 2,4,6-Tribromophenol	10.79	330	17315	6.95	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	99020	6.98	ng	0.00
78) Terphenyl-d14	13.09	244	64328	6.71	ng	0.00

Target Compounds

						Qvalue
48) Acenaphthylene	9.86	152	473466	20.647	ng	99
70) Phenanthrene	11.52	178	51436	2.754	ng	# 90
71) Anthracene	11.57	178	41039	2.104	ng	# 87
74) Fluoranthene	12.72	202	70896	3.571	ng	84
77) Pyrene	12.96	202	145832	9.158	ng	# 94
80) Benzo(a)anthracene	14.16	228	41231	2.975	ng	# 50
82) Chrysene	14.20	228	68402m	5.039	ng	
85) Indeno(1,2,3-cd)pyrene	17.33	276	65960	4.689	ng	# 89
87) Benzo(b)fluoranthene	15.29	252	93258m	5.954	ng	
88) Benzo(k)fluoranthene	15.30	252	49498m	3.355	ng	
89) Benzo(a)pyrene	15.68	252	81845	5.638	ng	# 88
91) Benzo(a,h,i)perylene	17.82	276	61195	4.552	ng	95

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

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