

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
 Data File : BF103417.D
 Acq On : 5 Mar 2018 19:10
 Operator : SJ/JU
 Sample : J1756-01 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-01-030118-A

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:23:10 PM

Quant Time: Mar 06 00:28:20 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 05 13:40:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	139574	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	574107	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	245978	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	386449	20.00	ng	0.00
75) Chrysene-d12	14.07	240	241205	20.00	ng	0.00
86) Perylene-d12	15.58	264	206787	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	401238	43.48	ng	0.00
7) Phenol-d6	6.51	99	523421	46.35	ng	-0.01
23) Nitrobenzene-d5	7.43	82	318853	31.72	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	64476	30.97	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	487400	28.24	ng	0.00
78) Terphenyl-d14	12.99	244	334976	33.38	ng	0.00
Target Compounds						Qvalue
70) Phenanthrene	11.43	178	246112	11.572	ng	98
71) Anthracene	11.49	178	77027	3.532	ng	99
74) Fluoranthene	12.63	202	351749	16.459	ng	98
77) Pyrene	12.86	202	312887	16.638	ng	99
80) Benzo(a)anthracene	14.06	228	123526	7.893	ng	98
82) Chrysene	14.09	228	119539	7.866	ng	98
85) Indeno(1,2,3-cd)pyrene	17.13	276	45621	3.792	ng	98
87) Benzo(b)fluoranthene	15.13	252	127503m	9.378	ng	
88) Benzo(k)fluoranthene	15.16	252	37570m	2.934	ng	
89) Benzo(a)pyrene	15.52	252	89212	7.348	ng	# 89
91) Benzo(g,h,i)perylene	17.60	276	44831	4.889	ng	# 91

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

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