

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
 Data File : BF103410.D
 Acq On : 5 Mar 2018 15:30
 Operator : SJ/JU
 Sample : J1751-11 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-030118-F

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:22:59 PM

Quant Time: Mar 05 17:05:38 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	135920	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	538186	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	216516	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	310835	20.00	ng	0.00
75) Chrysene-d12	14.07	240	245061	20.00	ng	0.00
86) Perylene-d12	15.59	264	173059	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	414762	46.15	ng	0.00
7) Phenol-d6	6.52	99	506221	46.03	ng	0.00
23) Nitrobenzene-d5	7.43	82	286865	30.44	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	58537	31.94	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	419459	27.50	ng	0.00
78) Terphenyl-d14	13.00	244	265240	26.02	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	51118	2.943	ng	Qvalue # 99
70) Phenanthrene	11.43	178	98841	5.778	ng	99
71) Anthracene	11.49	178	36854	2.101	ng	99
74) Fluoranthene	12.63	202	174808	10.169	ng	99
77) Pyrene	12.87	202	170853	8.942	ng	100
80) Benzo(a)anthracene	14.06	228	87034	5.474	ng	98
82) Chrysene	14.09	228	74931	4.853	ng	97
85) Indeno(1,2,3-cd)pyrene	17.13	276	24975	2.043	ng	96
87) Benzo(b)fluoranthene	15.14	252	82957m	7.291	ng	
88) Benzo(k)fluoranthene	15.16	252	24947m	2.328	ng	
89) Benzo(a)pyrene	15.52	252	53409	5.256	ng	# 86
91) Benzo(a,h,i)perylene	17.60	276	26079	3.399	ng	94

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

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