

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
 Data File : BF103401.D
 Acq On : 5 Mar 2018 11:31
 Operator : SJ/JU
 Sample : J1724-05
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-10

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:22:35 PM

Quant Time: Mar 05 13:47:39 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	152624	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	635400	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	279105	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	460506	20.00	ng	0.00
75) Chrysene-d12	14.06	240	278743	20.00	ng	0.00
86) Perylene-d12	15.58	264	281359	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	1082800	107.30	ng	0.00
7) Phenol-d6	6.52	99	1257348	101.82	ng	0.00
23) Nitrobenzene-d5	7.44	82	790375	71.04	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	207815	87.97	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1078469	62.39	ng	0.00
78) Terphenyl-d14	12.99	244	834098	71.93	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.53	94	107288	7.484	ng	87
49) Dimethylphthalate	9.62	163	102220	4.565	ng	99
70) Phenanthrene	11.43	178	227152	8.963	ng	99
71) Anthracene	11.49	178	65952	2.538	ng	99
74) Fluoranthene	12.63	202	436497	17.140	ng	99
77) Pyrene	12.86	202	409978	18.865	ng	99
80) Benzo(a)anthracene	14.05	228	188517	10.423	ng	98
82) Chrysene	14.09	228	173384	9.873	ng	97
85) Indeno(1,2,3-cd)pyrene	17.12	276	102362	7.363	ng	97
87) Benzo(b)fluoranthene	15.13	252	243575m	13.167	ng	
88) Benzo(k)fluoranthene	15.16	252	77595m	4.454	ng	
89) Benzo(a)pyrene	15.52	252	170518	10.322	ng	# 94
90) Dibenzo(a,h)anthracene	17.12	278	28919	2.265	ng	# 81
91) Benzo(g,h,i)perylene	17.59	276	105465	8.454	ng	95

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

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