

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
 Data File : BF103402.D
 Acq On : 5 Mar 2018 11:57
 Operator : SJ/JU
 Sample : J1724-11
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-19

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:22:38 PM

Quant Time: Mar 05 13:50:41 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	140958	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	590853	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	264900	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	433217	20.00	ng	0.00
75) Chrysene-d12	14.06	240	260173	20.00	ng	0.00
86) Perylene-d12	15.58	264	260578	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1004176	107.74	ng	0.00
7) Phenol-d6	6.52	99	1186776	104.06	ng	0.00
23) Nitrobenzene-d5	7.44	82	744917	72.00	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	193005	86.08	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	939333	56.20	ng	0.00
78) Terphenyl-d14	12.99	244	723504	66.85	ng	0.00
Target Compounds						
10) Phenol	6.53	94	32316	2.441	ng	Qvalue # 65
49) Dimethylphthalate	9.62	163	82299	3.873	ng	99
70) Phenanthrene	11.43	178	359786	15.091	ng	99
71) Anthracene	11.48	178	102242	4.182	ng	99
74) Fluoranthene	12.63	202	640313	26.727	ng	99
77) Pyrene	12.86	202	564860	27.847	ng	99
80) Benzo(a)anthracene	14.05	228	274551	16.264	ng	99
82) Chrysene	14.09	228	259449	15.829	ng	99
85) Indeno(1,2,3-cd)pyrene	17.12	276	110858	8.543	ng	97
87) Benzo(b)fluoranthene	15.13	252	322923m	18.848	ng	
88) Benzo(k)fluoranthene	15.16	252	98442m	6.102	ng	
89) Benzo(a)pyrene	15.52	252	212047	13.860	ng	# 92
90) Dibenzo(a,h)anthracene	17.12	278	31632	2.676	ng	# 74
91) Benzo(g,h,i)perylene	17.59	276	102210	8.846	ng	94

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

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