

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
 Data File : BF103427.D
 Acq On : 5 Mar 2018 23:35
 Operator : SJ/JU
 Sample : J1723-17
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-39

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:23:43 PM

Quant Time: Mar 06 01:22:58 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	126709	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	506023	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	201386	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	296763	20.00	ng	0.00
75) Chrysene-d12	14.07	240	176899	20.00	ng	0.00
86) Perylene-d12	15.59	264	131745	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	827122	98.73	ng	0.00
7) Phenol-d6	6.52	99	993578	96.92	ng	0.00
23) Nitrobenzene-d5	7.44	82	586868	66.23	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	119154	69.90	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	802046	64.76	ng	0.00
78) Terphenyl-d14	13.00	244	548528	74.54	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.18	128	111582	4.504	ng	99
37) 2-Methylnaphthalene	8.87	142	50077	3.128	ng	96
51) Acenaphthene	9.95	154	75937	6.340	ng	98
54) Dibenzofuran	10.12	168	90352	5.495	ng	96
57) Fluorene	10.47	166	93173	6.652	ng	98
70) Phenanthrene	11.44	178	677789	41.501	ng	99
71) Anthracene	11.49	178	221651	13.235	ng	99
72) Carbazole	11.65	167	65763	3.943	ng	98
74) Fluoranthene	12.63	202	760497	46.339	ng	99
77) Pyrene	12.87	202	695218	50.407	ng	100
80) Benzo(a)anthracene	14.06	228	300696	26.198	ng	99
82) Chrysene	14.10	228	266538	23.916	ng	98
85) Indeno(1,2,3-cd)pyrene	17.13	276	83405	9.453	ng	99
87) Benzo(b)fluoranthene	15.14	252	240295m	27.741	ng	
88) Benzo(k)fluoranthene	15.16	252	80913m	9.920	ng	
89) Benzo(a)pyrene	15.52	252	165525	21.399	ng	# 92
90) Dibenzo(a,h)anthracene	17.13	278	24422	4.086	ng	# 78
91) Benzo(g,h,i)perylene	17.60	276	77665	13.295	ng	# 89

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

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