

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
 Data File : BF103450.D
 Acq On : 6 Mar 2018 12:43
 Operator : SJ/JU
 Sample : J1723-17
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-39

Manual Integrations
 APPROVED

Sohil
 3/7/2018 12:14:45 PM

Quant Time: Mar 06 14:54:08 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 06 11:52:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	121338	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	514335	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	221024	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	331480	20.00	ng	0.00
75) Chrysene-d12	14.07	240	212875	20.00	ng	0.00
86) Perylene-d12	15.58	264	224117	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	790331	98.51	ng	0.00
7) Phenol-d6	6.51	99	978236	99.65	ng	0.00
23) Nitrobenzene-d5	7.44	82	609542	67.68	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	145494	77.77	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	849755	62.01	ng	0.00
78) Terphenyl-d14	12.99	244	556120	62.80	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.17	128	105511	4.190	ng	98
37) 2-Methylnaphthalene	8.87	142	51761	3.181	ng	100
51) Acenaphthene	9.95	154	82490	6.275	ng	99
54) Dibenzofuran	10.12	168	99382	5.507	ng	98
57) Fluorene	10.46	166	107725	7.008	ng	98
70) Phenanthrene	11.43	178	778977	42.702	ng	99
71) Anthracene	11.49	178	242953	12.988	ng	99
72) Carbazole	11.64	167	70247	3.771	ng	99
74) Fluoranthene	12.63	202	785799	42.866	ng	98
77) Pyrene	12.86	202	697022	41.997	ng	99
80) Benzo(a)anthracene	14.06	228	369059	26.720	ng	99
82) Chrysene	14.09	228	304345	22.693	ng	98
85) Indeno(1,2,3-cd)pyrene	17.13	276	129823	12.227	ng	97
87) Benzo(b)fluoranthene	15.14	252	408168m	27.700	ng	
88) Benzo(k)fluoranthene	15.16	252	107090m	7.718	ng	
89) Benzo(a)pyrene	15.52	252	274418	20.854	ng	# 91
90) Dibenzo(a,h)anthracene	17.12	278	35000	3.442	ng	# 82
91) Benzo(g,h,i)perylene	17.60	276	114496	11.522	ng	93

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

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