

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
 Data File : BF103452.D
 Acq On : 6 Mar 2018 13:37
 Operator : SJ/JU
 Sample : J1723-19
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-45

Manual Integrations
 APPROVED

Sohil
 3/7/2018 12:14:49 PM

Quant Time: Mar 06 14:59:58 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	120103	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	505393	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	212064	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	323301	20.00	ng	0.00
75) Chrysene-d12	14.07	240	212234	20.00	ng	0.00
86) Perylene-d12	15.58	264	179788	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	783894	98.71	ng	0.00
7) Phenol-d6	6.51	99	956007	98.38	ng	0.00
23) Nitrobenzene-d5	7.44	82	586000	66.22	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	126209	70.31	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	813847	61.87	ng	0.00
78) Terphenyl-d14	12.99	244	517291	58.59	ng	0.00
Target Compounds						Qvalue
51) Acenaphthene	9.95	154	26720	2.118	ng	98
57) Fluorene	10.46	166	31213	2.116	ng	99
70) Phenanthrene	11.43	178	282762	15.892	ng	98
71) Anthracene	11.49	178	96796	5.305	ng	99
72) Carbazole	11.65	167	37265	2.051	ng	94
74) Fluoranthene	12.63	202	385016	21.534	ng	98
77) Pyrene	12.86	202	382113	23.092	ng	99
80) Benzo(a)anthracene	14.06	228	184562	13.403	ng	99
82) Chrysene	14.09	228	173881	13.005	ng	97
85) Indeno(1,2,3-cd)pyrene	17.13	276	58350	5.512	ng	96
87) Benzo(b)fluoranthene	15.13	252	182309m	15.423	ng	
88) Benzo(k)fluoranthene	15.16	252	68013m	6.110	ng	
89) Benzo(a)pyrene	15.52	252	123256	11.676	ng	# 93
91) Benzo(a,h,i)perylene	17.60	276	56568	7.096	ng	93

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

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