

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
 Data File : BF103406.D
 Acq On : 5 Mar 2018 13:43
 Operator : SJ/JU
 Sample : J1751-07 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-030118-D

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:22:48 PM

Quant Time: Mar 05 15:21:49 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	137015	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	581027	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	255887	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	404779	20.00	ng	0.00
75) Chrysene-d12	14.07	240	252797	20.00	ng	0.00
86) Perylene-d12	15.58	264	228631	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	437654	48.31	ng	0.00
7) Phenol-d6	6.51	99	546233	49.28	ng	-0.01
23) Nitrobenzene-d5	7.43	82	322545	31.70	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	81022	37.41	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	476792	26.27	ng	0.00
78) Terphenyl-d14	12.99	244	336893	32.04	ng	0.00

Target Compounds

						Qvalue
37) 2-Methylnaphthalene	8.87	142	46762	2.544	ng	98
49) Dimethylphthalate	9.62	163	149940	7.304	ng	99
51) Acenaphthene	9.95	154	125378	8.238	ng	98
54) Dibenzofuran	10.12	168	178922	8.564	ng	98
57) Fluorene	10.46	166	242392	13.620	ng	100
70) Phenanthrene	11.44	178	1137860	51.080	ng	99
71) Anthracene	11.49	178	481987	21.100	ng	100
72) Carbazole	11.64	167	95301	4.189	ng	99
74) Fluoranthene	12.63	202	1167123	52.138	ng	98
77) Pyrene	12.86	202	999290	50.701	ng	100
80) Benzo(a)anthracene	14.06	228	513932	31.333	ng	99
82) Chrysene	14.09	228	396517	24.897	ng	98
85) Indeno(1,2,3-cd)pyrene	17.13	276	139810	11.089	ng	97
87) Benzo(b)fluoranthene	15.14	252	496543m	33.032	ng	
88) Benzo(k)fluoranthene	15.16	252	128223m	9.058	ng	
89) Benzo(a)pyrene	15.52	252	316817	23.601	ng	96
90) Dibenzo(a,h)anthracene	17.13	278	38581	3.719	ng	# 77
91) Benzo(g,h,i)perylene	17.60	276	128180	12.644	ng	97

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

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