

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111618\
 Data File : BF110838.D
 Acq On : 17 Nov 2018 1:59
 Operator : JU/SJ
 Sample : J5970-03 2X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-028-B

Manual Integrations
 APPROVED

Sohil
 11/19/2018 11:56:13 AM

Quant Time: Nov 17 03:50:57 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	64730	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	268596	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	118544	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	182835	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	156481	20.00	ng	-0.01
87) Perylene-d12	15.66	264	105523	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	192960	48.42	ng	-0.01
7) Phenol-d6	6.56	99	254897	50.02	ng	-0.01
23) Nitrobenzene-d5	7.50	82	157488	33.77	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	43579	36.48	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	305893	36.85	ng	-0.01
79) Terphenyl-d14	13.06	244	225993	27.05	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.57	94	12469	2.110	ng	86
50) Dimethylphthalate	9.69	163	51402	5.811	ng	96
52) Acenaphthene	10.01	154	18881	2.604	ng	96
55) Dibenzofuran	10.19	168	26972	2.543	ng	95
58) Fluorene	10.53	166	49656	6.094	ng	99
71) Phenanthrene	11.50	178	383744	39.176	ng	98
72) Anthracene	11.55	178	117113	11.852	ng	99
73) Carbazole	11.71	167	21302	2.245	ng	100
75) Fluoranthene	12.70	202	468381	47.694	ng	97
78) Pyrene	12.93	202	367129	30.470	ng	98
81) Benzo(a)anthracene	14.12	228	189500	20.261	ng	98
83) Chrysene	14.15	228	152136	17.073	ng	97
86) Indeno(1,2,3-cd)pyrene	17.23	276	36337	5.683	ng	95
88) Benzo(b)fluoranthene	15.20	252	141216m	22.370	ng	
89) Benzo(k)fluoranthene	15.23	252	40556m	6.502	ng	
90) Benzo(a)pyrene	15.59	252	83529	14.563	ng	98
91) Dibenzo(a,h)anthracene	17.23	278	10933m	2.252	ng	
92) Benzo(g,h,i)perylene	17.71	276	34359	7.135	ng	96

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

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