

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112918\
 Data File : BF111096.D
 Acq On : 29 Nov 2018 17:52
 Operator : JU/SJ
 Sample : J5767-19 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-112818-B

Manual Integrations
 APPROVED

mohammad
 11/30/2018 3:49:41 PM

Quant Time: Nov 30 03:00:40 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 29 10:35:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	67907	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	284559	20.00	ng	0.00
39) Acenaphthene-d10	9.97	164	126906	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	190976	20.00	ng	0.00
76) Chrysene-d12	14.13	240	154874	20.00	ng	0.00
87) Perylene-d12	15.66	264	120544	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	194600	47.52	ng	0.00
7) Phenol-d6	6.56	99	263071	49.30	ng	-0.01
23) Nitrobenzene-d5	7.50	82	162886	32.76	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	49575	39.33	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	272698	31.13	ng	0.00
79) Terphenyl-d14	13.04	244	185210	23.28	ng	-0.01
Target Compounds						
10) Phenol	6.57	94	21047	3.502	ng	82
50) Dimethylphthalate	9.69	163	23764	2.553	ng	99
71) Phenanthrene	11.49	178	44793	4.429	ng	97
75) Fluoranthene	12.69	202	99423	9.564	ng	97
78) Pyrene	12.92	202	96422	8.524	ng	98
81) Benzo(a)anthracene	14.12	228	62937	6.883	ng	97
83) Chrysene	14.15	228	52918	5.879	ng	99
86) Indeno(1,2,3-cd)pyrene	17.26	276	26133	4.045	ng	97
88) Benzo(b)fluoranthene	15.20	252	69156m	9.766	ng	
89) Benzo(k)fluoranthene	15.23	252	26471m	3.632	ng	
90) Benzo(a)pyrene	15.60	252	47879	7.264	ng	# 97
92) Benzo(a,h,i)perylene	17.74	276	23780	4.535	ng	93

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

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