

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102518\
 Data File : BF110183.D
 Acq On : 26 Oct 2018 1:41
 Operator : JU/SJ
 Sample : J5624-09 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PA-SB-01A

Manual Integrations
 APPROVED

Sohil
 10/26/2018 3:49:32 PM

Quant Time: Oct 26 07:18:08 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 23 15:06:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	136752	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	524185	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	212517	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	341606	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	241189	20.00	ng	-0.01
87) Perylene-d12	15.68	264	186159	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	330639	39.37	ng	-0.02
7) Phenol-d6	6.58	99	416828	38.81	ng	-0.02
23) Nitrobenzene-d5	7.53	82	233302	26.40	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	67120	31.88	ng	-0.02
45) 2-Fluorobiphenyl	9.32	172	433720	32.05	ng	-0.02
79) Terphenyl-d14	13.09	244	322777	27.83	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	56458	4.917	ng	Qvalue # 62
50) Dimethylphthalate	9.72	163	59224	3.775	ng	98
71) Phenanthrene	11.53	178	83636	4.679	ng	97
75) Fluoranthene	12.72	202	123996	6.835	ng	98
78) Pyrene	12.95	202	100613	5.819	ng	97
81) Benzo(a)anthracene	14.14	228	46913	3.280	ng	94
83) Chrysene	14.17	228	43226	3.182	ng	98
86) Indeno(1,2,3-cd)pyrene	17.23	276	24787	2.199	ng	96
88) Benzo(b)fluoranthene	15.22	252	53145m	4.944	ng	
90) Benzo(a)pyrene	15.61	252	29488	2.981	ng	# 92
92) Benzo(g,h,i)perylene	17.72	276	23436	2.786	ng	# 91

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

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