

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102618\
 Data File : BF110210.D
 Acq On : 26 Oct 2018 15:29
 Operator : JU/SJ
 Sample : J5624-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PA-SB-01A

Manual Integrations
 APPROVED

Sohil
 10/29/2018 11:56:39 AM

Quant Time: Oct 26 23:51:40 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	148589	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	575860	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	230749	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	345280	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	255184	20.00	ng	-0.01
87) Perylene-d12	15.68	264	175288	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	701021	76.83	ng	0.00
7) Phenol-d6	6.59	99	887713	76.06	ng	-0.01
23) Nitrobenzene-d5	7.53	82	549147	56.56	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	142470	62.33	ng	-0.01
45) 2-Fluorobiphenyl	9.33	172	925092	62.95	ng	-0.01
79) Terphenyl-d14	13.09	244	630323	51.37	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	126641	10.151	ng	Qvalue # 64
50) Dimethylphthalate	9.72	163	139935	8.215	ng	99
71) Phenanthrene	11.53	178	174515	9.660	ng	98
75) Fluoranthene	12.72	202	248563	13.556	ng	99
78) Pyrene	12.95	202	199385	10.899	ng	98
81) Benzo(a)anthracene	14.14	228	95660	6.321	ng	95
83) Chrysene	14.17	228	104123	7.244	ng	96
84) Bis(2-ethylhexyl)phthalate	14.11	149	22350	2.082	ng	# 93
86) Indeno(1,2,3-cd)pyrene	17.24	276	43032	3.609	ng	96
88) Benzo(b)fluoranthene	15.23	252	121586m	12.012	ng	
89) Benzo(k)fluoranthene	15.25	252	32311m	3.247	ng	
90) Benzo(a)pyrene	15.61	252	58745	6.307	ng	# 92
92) Benzo(a,h,i)perylene	17.72	276	41835	5.283	ng	# 90

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

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