

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102418\
 Data File : BF110112.D
 Acq On : 24 Oct 2018 13:57
 Operator : JU/SJ
 Sample : J5610-05 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3(0-2)

Manual Integrations
 APPROVED

Sohil
 10/25/2018 11:14:33 AM

Quant Time: Oct 25 03:59:58 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	175299	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	660735	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	274150	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	440408	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	350973	20.00	ng	-0.01
87) Perylene-d12	15.68	264	254616	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	504058	46.82	ng	-0.02
7) Phenol-d6	6.58	99	621826	45.16	ng	-0.02
23) Nitrobenzene-d5	7.53	82	353773	31.76	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	100736	37.09	ng	-0.02
45) 2-Fluorobiphenyl	9.32	172	554296	31.75	ng	-0.02
79) Terphenyl-d14	13.09	244	436779	25.88	ng	-0.02
Target Compounds						
10) Phenol	6.59	94	55008	3.737	ng	Qvalue # 63
50) Dimethylphthalate	9.72	163	97769	4.831	ng	98
71) Phenanthrene	11.53	178	156317	6.783	ng	98
72) Anthracene	11.57	178	51772	2.241	ng	98
75) Fluoranthene	12.72	202	323783	13.844	ng	99
78) Pyrene	12.95	202	296320	11.777	ng	97
81) Benzo(a)anthracene	14.14	228	170487	8.191	ng	97
83) Chrysene	14.17	228	145902	7.380	ng	97
86) Indeno(1,2,3-cd)pyrene	17.24	276	65704	4.006	ng	99
88) Benzo(b)fluoranthene	15.22	252	162952m	11.083	ng	
89) Benzo(k)fluoranthene	15.25	252	56871m	3.934	ng	
90) Benzo(a)pyrene	15.61	252	104063	7.692	ng	# 94
92) Benzo(a,h,i)perylene	17.72	276	60395	5.250	ng	# 93

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

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