

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102618\
 Data File : BF110213.D
 Acq On : 26 Oct 2018 16:50
 Operator : JU/SJ
 Sample : J5624-06
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PA-SB-03B

Manual Integrations
 APPROVED

Sohil
 10/29/2018 11:56:40 AM

Quant Time: Oct 27 00:03:28 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	120677	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	472053	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	199814	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	310770	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	232163	20.00	ng	-0.01
87) Perylene-d12	15.68	264	164171	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	733296	98.95	ng	0.00
7) Phenol-d6	6.59	99	931766	98.30	ng	-0.01
23) Nitrobenzene-d5	7.53	82	574133	72.14	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	172530	87.17	ng	-0.01
45) 2-Fluorobiphenyl	9.33	172	918874	72.21	ng	-0.01
79) Terphenyl-d14	13.09	244	668889	59.92	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	130302	12.860	ng	Qvalue # 63
50) Dimethylphthalate	9.72	163	157544	10.681	ng	99
71) Phenanthrene	11.53	178	154495	9.501	ng	99
75) Fluoranthene	12.72	202	158523	9.606	ng	98
78) Pyrene	12.95	202	146255	8.787	ng	98
81) Benzo(a)anthracene	14.14	228	65837	4.782	ng	99
83) Chrysene	14.17	228	58409	4.467	ng	97
86) Indeno(1,2,3-cd)pyrene	17.24	276	25092	2.313	ng	96
88) Benzo(b)fluoranthene	15.22	252	59134m	6.238	ng	
90) Benzo(a)pyrene	15.61	252	36153	4.144	ng	# 90
92) Benzo(g,h,i)perylene	17.72	276	22073	2.976	ng	# 93

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

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