

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102418\
 Data File : BF110104.D
 Acq On : 24 Oct 2018 10:19
 Operator : JU/SJ
 Sample : J5610-08
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-4(8-10)

Manual Integrations
 APPROVED

Sohil
 10/25/2018 11:14:01 AM

Quant Time: Oct 25 02:35:00 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	6.96	152	156955	20.00	ng	-0.02
21) Naphthalene-d8	8.25	136	602760	20.00	ng	-0.02
39) Acenaphthene-d10	10.00	164	261404	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	401601	20.00	ng	-0.02
76) Chrysene-d12	14.14	240	298211	20.00	ng	-0.02
87) Perylene-d12	15.66	264	314831	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	866964	89.95	ng	-0.01
7) Phenol-d6	6.59	99	1054503	85.54	ng	-0.02
23) Nitrobenzene-d5	7.53	82	664294	65.37	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	204396	78.94	ng	-0.02
45) 2-Fluorobiphenyl	9.32	172	1001117	60.14	ng	-0.02
79) Terphenyl-d14	13.08	244	643942	44.91	ng	-0.02
Target Compounds						
10) Phenol	6.59	94	94766	7.191	ng	Qvalue # 73
50) Dimethylphthalate	9.71	163	173165	8.974	ng	98
71) Phenanthrene	11.52	178	104218	4.960	ng	97
75) Fluoranthene	12.71	202	104852	4.917	ng	100
78) Pyrene	12.94	202	118391	5.538	ng	96
81) Benzo(a)anthracene	14.13	228	63725	3.603	ng	97
83) Chrysene	14.17	228	61229	3.645	ng	96
84) Bis(2-ethylhexyl)phthalate	14.10	149	45560	3.632	ng	96
86) Indeno(1,2,3-cd)pyrene	17.21	276	35456	2.544	ng	# 93
88) Benzo(b)fluoranthene	15.21	252	70990m	3.905	ng	
90) Benzo(a)pyrene	15.59	252	53752	3.213	ng	# 94
92) Benzo(a,h,i)perylene	17.68	276	34131	2.400	ng	# 88

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

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