

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102518\
 Data File : BF110176.D
 Acq On : 25 Oct 2018 22:29
 Operator : JU/SJ
 Sample : J5625-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DW-06

Manual Integrations
 APPROVED

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

Quant Time: Oct 26 05:28:28 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	6.96	152	162769	20.00	ng	-0.02
21) Naphthalene-d8	8.25	136	638072	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	266524	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	400989	20.00	ng	-0.02
76) Chrysene-d12	14.14	240	293532	20.00	ng	-0.02
87) Perylene-d12	15.67	264	254843	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	930784	93.12	ng	0.00
7) Phenol-d6	6.59	99	1180757	92.36	ng	-0.01
23) Nitrobenzene-d5	7.53	82	754447	70.13	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	209504	79.35	ng	-0.02
45) 2-Fluorobiphenyl	9.32	172	1178905	69.46	ng	-0.02
79) Terphenyl-d14	13.09	244	765643	54.25	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	136459	9.985	ng	Qvalue # 62
31) Naphthalene	8.27	128	60917	2.071	ng	99
37) 2-Methylnaphthalene	8.96	142	39256	2.018	ng	98
50) Dimethylphthalate	9.72	163	215241	10.940	ng	100
71) Phenanthrene	11.52	178	128575	6.128	ng	99
72) Anthracene	11.57	178	46966	2.233	ng	98
75) Fluoranthene	12.72	202	239227	11.235	ng	100
78) Pyrene	12.94	202	292787	13.913	ng	98
80) Butylbenzylphthalate	13.55	149	67980	7.443	ng	99
81) Benzo(a)anthracene	14.14	228	164802m	9.468	ng	
83) Chrysene	14.17	228	173447	10.491	ng	98
84) Bis(2-ethylhexyl)phthalate	14.10	149	29351	2.377	ng	# 95
86) Indeno(1,2,3-cd)pyrene	17.23	276	74599	5.439	ng	100
88) Benzo(b)fluoranthene	15.22	252	225914m	15.351	ng	
89) Benzo(k)fluoranthene	15.24	252	53633m	3.707	ng	
90) Benzo(a)pyrene	15.60	252	129591	9.570	ng	97
92) Benzo(a,h,i)perylene	17.70	276	75299	6.540	ng	# 94

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

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