

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111218\
 Data File : BF110691.D
 Acq On : 12 Nov 2018 22:54
 Operator : JU/SJ
 Sample : J5907-20 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2-S

Manual Integrations
 APPROVED

Sohil
 11/13/2018 2:22:55 PM

Quant Time: Nov 13 06:38:30 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 09 15:13:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	93778	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	399254	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	180832	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	294712	20.00	ng	0.00
76) Chrysene-d12	14.13	240	185446	20.00	ng	-0.01
87) Perylene-d12	15.66	264	167057	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	296173	51.30	ng	0.00
7) Phenol-d6	6.56	99	387145	52.44	ng	-0.01
23) Nitrobenzene-d5	7.50	82	243121	35.07	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	64463	35.38	ng	-0.01
45) 2-Fluorobiphenyl	9.30	172	459960	36.33	ng	0.00
79) Terphenyl-d14	13.06	244	319495	32.26	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.57	94	21524	2.514	ng	80
50) Dimethylphthalate	9.69	163	64404	4.773	ng	100
71) Phenanthrene	11.50	178	243842	15.444	ng	99
72) Anthracene	11.56	178	55978	3.515	ng	97
75) Fluoranthene	12.70	202	272609	17.221	ng	97
78) Pyrene	12.93	202	242545	16.986	ng	98
81) Benzo(a)anthracene	14.12	228	113341	10.225	ng	99
83) Chrysene	14.15	228	102134	9.672	ng	97
86) Indeno(1,2,3-cd)pyrene	17.22	276	42925	5.665	ng	95
88) Benzo(b)fluoranthene	15.20	252	120865m	12.094	ng	
89) Benzo(k)fluoranthene	15.23	252	32595m	3.301	ng	
90) Benzo(a)pyrene	15.59	252	79530	8.759	ng	97
92) Benzo(a,h,i)perylene	17.70	276	40723	5.341	ng	# 90

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

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