

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
 Data File : BF111621.D
 Acq On : 20 Dec 2018 5:31
 Operator : JU/SJ
 Sample : J6340-07
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-05A

Manual Integrations
 APPROVED

Sohil
 12/20/2018 12:42:46 PM

Quant Time: Dec 20 08:12:34 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	198445	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	688919	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	318302	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	577286	20.00	ng	0.00
76) Chrysene-d12	14.05	240	508941	20.00	ng	0.00
87) Perylene-d12	15.52	264	356383	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	1502186	129.03	ng	0.01
7) Phenol-d6	6.53	99	1893159	127.77	ng	0.00
23) Nitrobenzene-d5	7.46	82	1169420	98.81	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	492458	130.94	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	2091347	95.77	ng	0.00
79) Terphenyl-d14	13.00	244	2069509	77.11	ng	0.00
Target Compounds						
10) Phenol	6.54	94	98843	5.913	ng	94
31) Naphthalene	8.20	128	97417	2.943	ng	95
37) 2-Methylnaphthalene	8.89	142	54127	2.513	ng	97
38) 1-Methylnaphthalene	8.99	142	42478	2.054	ng	100
50) Dimethylphthalate	9.64	163	248626	10.681	ng	97
75) Fluoranthene	12.63	202	533349	17.203	ng	95
78) Pyrene	12.86	202	445946	12.408	ng	96
81) Benzo(a)anthracene	14.04	228	187732	6.463	ng	96
83) Chrysene	14.07	228	293084	10.267	ng	98
86) Indeno(1,2,3-cd)pyrene	17.00	276	62255	2.565	ng	# 85
88) Benzo(b)fluoranthene	15.09	252	278354m	13.364	ng	
89) Benzo(k)fluoranthene	15.12	252	77234m	3.895	ng	
90) Benzo(a)pyrene	15.46	252	90751	4.796	ng	94
92) Benzo(a,h,i)perylene	17.44	276	56195m	3.473	ng	

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

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