

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121318\
 Data File : BF111390.D
 Acq On : 14 Dec 2018 00:37
 Operator : JU/SJ
 Sample : J6210-07
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C-1-112918

Manual Integrations
 APPROVED

Sohil
 12/14/2018 12:53:49 PM

Quant Time: Dec 14 04:00:45 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 13:16:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	250739	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	1136012	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	450185	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	648738	20.00	ng	0.00
76) Chrysene-d12	13.95	240	453320	20.00	ng	0.00
87) Perylene-d12	15.39	264	348416	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	1300976	87.74	ng	0.02
7) Phenol-d6	6.45	99	1718860	85.98	ng	0.01
23) Nitrobenzene-d5	7.35	82	1044494	51.82	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	280613	70.71	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1886033	67.26	ng	0.00
79) Terphenyl-d14	12.90	244	1282124	62.25	ng	0.00
Target Compounds						
10) Phenol	6.45	94	122629	5.415	ng	97
50) Dimethylphthalate	9.55	163	424989	13.128	ng	98
71) Phenanthrene	11.33	178	93424	2.786	ng	# 90
75) Fluoranthene	12.52	202	206212	6.300	ng	99
78) Pyrene	12.75	202	217130	6.206	ng	94
81) Benzo(a)anthracene	13.94	228	97657	3.850	ng	94
83) Chrysene	13.97	228	116053	4.485	ng	97
88) Benzo(b)fluoranthene	14.97	252	137082m	6.919	ng	
90) Benzo(a)pyrene	15.32	252	83517	4.527	ng	# 95
92) Benzo(g,h,i)perylene	17.22	276	56478	3.681	ng	# 92

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

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