

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121218\
 Data File : BF111347.D
 Acq On : 13 Dec 2018 3:02
 Operator : JU/SJ
 Sample : J6214-49
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-7(0-2)

Manual Integrations
 APPROVED

Sohil
 12/13/2018 6:07:01 PM

Quant Time: Dec 13 06:05:52 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	6.80	152	226329	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	810826	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	367398	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	631297	20.00	ng	0.00
76) Chrysene-d12	13.95	240	397810	20.00	ng	0.00
87) Perylene-d12	15.39	264	326792	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1379101	103.04	ng	0.00
7) Phenol-d6	6.43	99	1710148	94.77	ng	0.00
23) Nitrobenzene-d5	7.35	82	921386	64.05	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	285668	88.21	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1701278	74.34	ng	0.00
79) Terphenyl-d14	12.90	244	1380873	76.40	ng	0.00
Target Compounds						
10) Phenol	6.44	94	79876	3.908	ng	93
50) Dimethylphthalate	9.55	163	156464	5.922	ng	96
71) Phenanthrene	11.34	178	72342	2.217	ng	94
75) Fluoranthene	12.53	202	221229	6.945	ng	96
78) Pyrene	12.75	202	211316	6.883	ng	95
81) Benzo(a)anthracene	13.94	228	81200	3.648	ng	98
83) Chrysene	13.97	228	77636	3.419	ng	97
88) Benzo(b)fluoranthene	14.98	252	86106m	4.634	ng	
90) Benzo(a)pyrene	15.33	252	57126	3.302	ng	# 85
92) Benzo(g,h,i)perylene	17.23	276	32424	2.253	ng	# 85

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

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