

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112318\
 Data File : BF110947.D
 Acq On : 23 Nov 2018 16:40
 Operator : JU/SJ
 Sample : J5999-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSSI-1115-S-AB-1

Manual Integrations
 APPROVED

MOHAMMAD
 11/26/2018 3:12:04 PM

Quant Time: Nov 23 21:19:53 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	58512	20.00	ng	-0.01
21) Naphthalene-d8	8.22	136	239138	20.00	ng	-0.01
39) Acenaphthene-d10	9.97	164	103223	20.00	ng	-0.02
64) Phenanthrene-d10	11.47	188	165404	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	138117	20.00	ng	-0.01
87) Perylene-d12	15.66	264	126677	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	354126	98.31	ng	0.00
7) Phenol-d6	6.56	99	447089	97.06	ng	-0.01
23) Nitrobenzene-d5	7.50	82	280716	67.60	ng	-0.02
42) 2,4,6-Tribromophenol	10.77	330	81546	78.40	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	356430	49.32	ng	-0.02
79) Terphenyl-d14	13.05	244	280603	38.05	ng	-0.02
Target Compounds						
10) Phenol	6.57	94	23198	4.343	ng	Qvalue # 74
50) Dimethylphthalate	9.69	163	31259	4.058	ng	96
71) Phenanthrene	11.50	178	21370	2.412	ng	98
75) Fluoranthene	12.70	202	56503	6.360	ng	96
78) Pyrene	12.93	202	51544	4.847	ng	98
81) Benzo(a)anthracene	14.12	228	31545	3.821	ng	98
83) Chrysene	14.15	228	30138	3.832	ng	99
86) Indeno(1,2,3-cd)pyrene	17.26	276	11940	2.116	ng	# 88
88) Benzo(b)fluoranthene	15.20	252	36659m	4.837	ng	
90) Benzo(a)pyrene	15.60	252	26055	3.784	ng	# 94

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

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