

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
 Data File : BF111714.D
 Acq On : 26 Dec 2018 14:23
 Operator : JU/SJ
 Sample : J6498-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSSI-1218-S-AB-6

Manual Integrations
 APPROVED

Sohil
 12/27/2018 8:22:19 AM

Quant Time: Dec 26 15:35:59 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 26 13:36:07 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	165825	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	614228	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	306275	20.00	ng	-0.01
64) Phenanthrene-d10	11.42	188	560296	20.00	ng	0.00
76) Chrysene-d12	14.06	240	495442	20.00	ng	-0.01
87) Perylene-d12	15.54	264	450644	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	810370	83.30	ng	0.00
7) Phenol-d6	6.54	99	1038053	83.84	ng	0.00
23) Nitrobenzene-d5	7.46	82	651233	61.71	ng	-0.01
42) 2,4,6-Tribromophenol	10.73	330	308484	85.24	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	1248712	59.43	ng	0.00
79) Terphenyl-d14	13.00	244	1322922	50.64	ng	0.00
Target Compounds						
10) Phenol	6.55	94	74157	5.308	ng	94
50) Dimethylphthalate	9.65	163	174430	7.788	ng	96
52) Acenaphthene	9.97	154	42934	2.328	ng	95
71) Phenanthrene	11.44	178	219484	7.386	ng	93
72) Anthracene	11.49	178	60372	2.024	ng	94
75) Fluoranthene	12.63	202	199415	6.627	ng	93
78) Pyrene	12.86	202	161890	4.627	ng	96
81) Benzo(a)anthracene	14.05	228	82838	2.930	ng	99
83) Chrysene	14.09	228	71348	2.567	ng	96
88) Benzo(b)fluoranthene	15.10	252	72743m	2.762	ng	
90) Benzo(a)pyrene	15.47	252	53030	2.216	ng	# 93

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

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