

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
 Data File : BF111599.D
 Acq On : 19 Dec 2018 19:25
 Operator : JU/SJ
 Sample : J6403-01RE
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1-A

Manual Integrations
 APPROVED

Sohil
 12/20/2018 12:42:01 PM

Quant Time: Dec 20 11:00:48 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	6.89	152	192591	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	703318	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	337462	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	601540	20.00	ng	0.00
76) Chrysene-d12	14.04	240	517146	20.00	ng	0.00
87) Perylene-d12	15.52	264	559493	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	875459	77.48	ng	0.00
7) Phenol-d6	6.52	99	1102375	76.66	ng	0.00
23) Nitrobenzene-d5	7.45	82	675502	55.91	ng	-0.01
42) 2,4,6-Tribromophenol	10.71	330	285085	71.50	ng	-0.01
45) 2-Fluorobiphenyl	9.25	172	1138023	49.15	ng	0.00
79) Terphenyl-d14	12.99	244	1002265	36.75	ng	0.00
Target Compounds						
10) Phenol	6.53	94	71304	4.395	ng	Qvalue 93
50) Dimethylphthalate	9.64	163	117703	4.770	ng	97
71) Phenanthrene	11.43	178	111822	3.505	ng	94
75) Fluoranthene	12.62	202	216906	6.714	ng	95
78) Pyrene	12.84	202	198533	5.436	ng	95
81) Benzo(a)anthracene	14.03	228	96105	3.256	ng	100
83) Chrysene	14.07	228	98807	3.406	ng	98
86) Indeno(1,2,3-cd)pyrene	16.98	276	60153	2.439	ng	# 84
88) Benzo(b)fluoranthene	15.09	252	133943m	4.096	ng	
90) Benzo(a)pyrene	15.45	252	92360m	3.109	ng	
92) Benzo(g,h,i)perylene	17.43	276	52028	2.048	ng	# 87

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

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