

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121718\
 Data File : BF111562.D
 Acq On : 18 Dec 2018 1:58
 Operator : JU/SJ
 Sample : J6403-45
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-6-B

Manual Integrations
 APPROVED

Sohil
 12/18/2018 12:49:53 PM

Quant Time: Dec 18 07:13:49 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	137451	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	499377	20.00	ng	-0.01
39) Acenaphthene-d10	9.82	164	226291	20.00	ng	-0.01
64) Phenanthrene-d10	11.31	188	337508	20.00	ng	0.00
76) Chrysene-d12	13.94	240	247714	20.00	ng	0.00
87) Perylene-d12	15.39	264	212123	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	922661	111.70	ng	0.01
7) Phenol-d6	6.44	99	1185923	107.94	ng	0.00
23) Nitrobenzene-d5	7.35	82	674927	80.48	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	198080	97.72	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1046791	71.45	ng	-0.01
79) Terphenyl-d14	12.89	244	684015	64.84	ng	0.00
Target Compounds						
10) Phenol	6.45	94	69850	5.707	ng	98
50) Dimethylphthalate	9.54	163	123414	7.573	ng	# 95
71) Phenanthrene	11.33	178	78072	4.489	ng	# 91
75) Fluoranthene	12.52	202	124008	7.289	ng	95
78) Pyrene	12.74	202	133245	7.629	ng	95
81) Benzo(a)anthracene	13.93	228	65735	4.880	ng	94
83) Chrysene	13.97	228	58843	4.147	ng	95
86) Indeno(1,2,3-cd)pyrene	16.79	276	25638	2.504	ng	94
88) Benzo(b)fluoranthene	14.97	252	71942m	5.817	ng	
89) Benzo(k)fluoranthene	14.99	252	24709m	2.091	ng	
90) Benzo(a)pyrene	15.32	252	53769	4.821	ng	97
92) Benzo(a,h,i)perylene	17.22	276	24948	2.889	ng	# 90

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

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