

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110218\
 Data File : BF110424.D
 Acq On : 2 Nov 2018 20:57
 Operator : JU/SJ
 Sample : J5767-05 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-110118-C

Manual Integrations
 APPROVED

Sohil
 11/5/2018 10:32:47 AM

Quant Time: Nov 03 00:40:39 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 16:06:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	87960	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	353314	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	141755	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	214044	20.00	ng	0.00
76) Chrysene-d12	14.13	240	171836	20.00	ng	0.00
87) Perylene-d12	15.66	264	127472	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	127748	23.65	ng	0.00
7) Phenol-d6	6.57	99	163375	23.65	ng	-0.01
23) Nitrobenzene-d5	7.52	82	95620	16.05	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	23083	16.44	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	171272	18.97	ng	0.00
79) Terphenyl-d14	13.06	244	114083	13.81	ng	-0.01
Target Compounds						Qvalue
50) Dimethylphthalate	9.70	163	25931	2.478	ng	98
71) Phenanthrene	11.51	178	31355	2.800	ng	97
75) Fluoranthene	12.70	202	55982	4.925	ng	99
78) Pyrene	12.93	202	53582	4.349	ng	97
81) Benzo(a)anthracene	14.12	228	32543	3.194	ng	96
83) Chrysene	14.16	228	29367	3.034	ng	96
88) Benzo(b)fluoranthene	15.20	252	31422m	4.269	ng	
90) Benzo(a)pyrene	15.59	252	21435	3.165	ng	# 91
92) Benzo(g,h,i)perylene	17.69	276	12710	2.207	ng	# 92

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

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