

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110718\
 Data File : BF110569.D
 Acq On : 7 Nov 2018 21:28
 Operator : JU/SJ
 Sample : J5813-10
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-J

Manual Integrations
 APPROVED

Sohil
 11/9/2018 1:35:19 PM

Quant Time: Nov 09 10:48:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 09 10:28:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	64209	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	224808	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	85246	20.00	ng	-0.01
64) Phenanthrene-d10	11.49	188	153430	20.00	ng	0.00
76) Chrysene-d12	14.14	240	156994	20.00	ng	0.00
87) Perylene-d12	15.67	264	103822	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	345038	87.51	ng	0.00
7) Phenol-d6	6.57	99	410717	81.44	ng	-0.01
23) Nitrobenzene-d5	7.52	82	245249	64.71	ng	-0.01
42) 2,4,6-Tribromophenol	10.79	330	70615	83.63	ng	-0.01
45) 2-Fluorobiphenyl	9.30	172	390248	71.88	ng	-0.01
79) Terphenyl-d14	13.07	244	362830	48.06	ng	-0.01
Target Compounds						
10) Phenol	6.59	94	24851	4.610	ng	Qvalue # 69
50) Dimethylphthalate	9.70	163	52200	8.295	ng	100
71) Phenanthrene	11.51	178	24513	3.053	ng	97
75) Fluoranthene	12.70	202	62154	7.628	ng	98
78) Pyrene	12.94	202	51142	4.544	ng	97
81) Benzo(a)anthracene	14.13	228	22379	2.404	ng	97
83) Chrysene	14.16	228	25697	2.906	ng	97
88) Benzo(b)fluoranthene	15.22	252	27585m	4.601	ng	
90) Benzo(a)pyrene	15.60	252	14383	2.607	ng	# 87
92) Benzo(g,h,i)perylene	17.73	276	10433	2.224	ng	# 93

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

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