

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110918\
 Data File : BF110617.D
 Acq On : 9 Nov 2018 10:31
 Operator : JU/SJ
 Sample : J5813-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-D

Manual Integrations
 APPROVED

Sohil
 11/12/2018 10:29:53 AM

Quant Time: Nov 09 15:53:18 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 09 15:13:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	80112	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	314650	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	129016	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	194701	20.00	ng	0.00
76) Chrysene-d12	14.13	240	160542	20.00	ng	-0.01
87) Perylene-d12	15.66	264	125295	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	537526	108.99	ng	0.00
7) Phenol-d6	6.57	99	670550	106.32	ng	0.00
23) Nitrobenzene-d5	7.51	82	426514	78.06	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	92833	71.41	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	711739	78.79	ng	0.00
79) Terphenyl-d14	13.06	244	470842	54.92	ng	0.00
Target Compounds						
10) Phenol	6.58	94	43012	5.881	ng	Qvalue # 71
50) Dimethylphthalate	9.69	163	99900	10.377	ng	99
71) Phenanthrene	11.50	178	21942	2.104	ng	97
75) Fluoranthene	12.70	202	46589	4.455	ng	98
78) Pyrene	12.93	202	40716	3.294	ng	99
83) Chrysene	14.16	228	20928	2.289	ng	95
88) Benzo(b)fluoranthene	15.20	252	27032m	3.606	ng	
90) Benzo(a)pyrene	15.59	252	15077	2.214	ng	# 84

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

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