

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\
 Data File : BF110720.D
 Acq On : 13 Nov 2018 15:47
 Operator : JU/SJ
 Sample : J5908-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SR-2-S

Manual Integrations
 APPROVED

Sohil
 11/14/2018 1:48:31 PM

Quant Time: Nov 13 16:49:09 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	76249	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	306993	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	136355	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	230157	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	153982	20.00	ng	-0.01
87) Perylene-d12	15.65	264	163339	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	460968	98.20	ng	0.00
7) Phenol-d6	6.57	99	569874	94.93	ng	0.00
23) Nitrobenzene-d5	7.51	82	353331	66.28	ng	0.00
42) 2,4,6-Tribromophenol	10.77	330	111534	81.18	ng	-0.01
45) 2-Fluorobiphenyl	9.30	172	587647	61.55	ng	0.00
79) Terphenyl-d14	13.06	244	451680	54.93	ng	-0.01
Target Compounds						
10) Phenol	6.58	94	25707	3.693	ng	Qvalue # 76
50) Dimethylphthalate	9.69	163	32127	3.158	ng	99
75) Fluoranthene	12.70	202	27169	2.198	ng	98
78) Pyrene	12.93	202	25049	2.113	ng	98
86) Indeno(1,2,3-cd)pyrene	17.23	276	13106	2.083	ng	97
88) Benzo(b)fluoranthene	15.20	252	25780m	2.638	ng	
90) Benzo(a)pyrene	15.59	252	19027	2.143	ng	# 93

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

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