

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\
 Data File : BF110730.D
 Acq On : 13 Nov 2018 20:21
 Operator : JU/SJ
 Sample : J5910-03 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-4-S

Manual Integrations
 APPROVED

Sohil
 11/14/2018 1:52:24 PM

Quant Time: Nov 14 05:56:09 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	91824	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	376950	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	154028	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	228490	20.00	ng	0.00
76) Chrysene-d12	14.13	240	186930	20.00	ng	-0.01
87) Perylene-d12	15.66	264	158322	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	269942	47.75	ng	0.00
7) Phenol-d6	6.56	99	340266	47.07	ng	-0.01
23) Nitrobenzene-d5	7.50	82	209484	32.00	ng	-0.01
42) 2,4,6-Tribromophenol	10.78	330	59132	38.10	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	373229	34.61	ng	0.00
79) Terphenyl-d14	13.06	244	239067	23.95	ng	-0.01
Target Compounds						
10) Phenol	6.58	94	17510	2.089	ng	Qvalue # 76
50) Dimethylphthalate	9.69	163	54359	4.730	ng	99
71) Phenanthrene	11.50	178	34179	2.792	ng	99
75) Fluoranthene	12.70	202	41079	3.347	ng	98
78) Pyrene	12.93	202	34537	2.400	ng	98
88) Benzo(b)fluoranthene	15.20	252	22181m	2.342	ng	

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

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