

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112718\
 Data File : BF111014.D
 Acq On : 27 Nov 2018 17:19
 Operator : JU/SJ
 Sample : J6065-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 USTS-1-2-3

Manual Integrations
 APPROVED

mohammad
 11/28/2018 4:42:24 PM

Quant Time: Nov 27 17:47:45 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 27 15:32:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	50557	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	181054	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	76438	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	140861	20.00	ng	0.00
76) Chrysene-d12	14.14	240	111298	20.00	ng	0.00
87) Perylene-d12	15.67	264	72310	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	382223	125.37	ng	0.01
7) Phenol-d6	6.57	99	453556	114.16	ng	0.00
23) Nitrobenzene-d5	7.50	82	263860	83.41	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	89520	117.92	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	419407	79.49	ng	0.00
79) Terphenyl-d14	13.06	244	416135	72.78	ng	0.00
Target Compounds						
10) Phenol	6.58	94	24903	5.566	ng	81
50) Dimethylphthalate	9.69	163	25376	4.527	ng	97
75) Fluoranthene	12.70	202	16271	2.122	ng	98
78) Pyrene	12.94	202	18438	2.268	ng	96
88) Benzo(b)fluoranthene	15.23	252	9771m	2.300	ng	

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

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