

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112818\
 Data File : BF111057.D
 Acq On : 28 Nov 2018 20:39
 Operator : JU/SJ
 Sample : J6079-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 J-1

Manual Integrations
 APPROVED

mohammad
 11/29/2018 1:58:35 PM

Quant Time: Nov 29 03:15:38 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	6.93	152	54981	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	219467	20.00	ng	0.00
39) Acenaphthene-d10	9.97	164	86612	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	135970	20.00	ng	0.00
76) Chrysene-d12	14.13	240	119813	20.00	ng	0.00
87) Perylene-d12	15.66	264	91963	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	360058	108.60	ng	0.00
7) Phenol-d6	6.56	99	448611	103.83	ng	0.00
23) Nitrobenzene-d5	7.50	82	278721	72.68	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	82715	96.16	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	489058	81.81	ng	0.00
79) Terphenyl-d14	13.05	244	339974	55.24	ng	0.00
Target Compounds						
10) Phenol	6.57	94	21118	4.340	ng	89
50) Dimethylphthalate	9.69	163	50563	7.960	ng	98
75) Fluoranthene	12.70	202	22875	3.091	ng	99
78) Pyrene	12.93	202	23174	2.648	ng	96
88) Benzo(b)fluoranthene	15.22	252	12512m	2.316	ng	

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

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