

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101819\
 Data File : BP000859.D
 Acq On : 18 Oct 2019 01:04
 Operator : JU
 Sample : K5401-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR200031-RT4221

Manual Integrations
 APPROVED

mohammad
 10/20/2019 8:16:38 PM

Quant Time: Oct 18 07:34:37 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 01 18:03:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	187232	20.00	ng	-0.02
21) Naphthalene-d8	8.03	136	668642	20.00	ng	-0.02
39) Acenaphthene-d10	9.80	164	368445	20.00	ng	-0.02
64) Phenanthrene-d10	11.29	188	678063	20.00	ng	-0.02
76) Chrysene-d12	13.97	240	587279	20.00	ng	-0.01
87) Perylene-d12	15.45	264	653937	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	955454	82.03	ng	-0.02
7) Phenol-d6	6.39	99	1192618	84.97	ng	-0.02
23) Nitrobenzene-d5	7.31	82	663351	60.59	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	360939	91.93	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	1306692	57.38	ng	-0.02
79) Terphenyl-d14	12.90	244	1575504	54.32	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.51	163	220501	8.613	ng	99
75) Fluoranthene	12.52	202	182734	4.778	ng	99
78) Pyrene	12.75	202	164763	4.068	ng	99
88) Benzo(b)fluoranthene	15.02	252	74567m	2.010	ng	

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

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