

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101819\
 Data File : BP000864.D
 Acq On : 18 Oct 2019 03:29
 Operator : JU
 Sample : K5401-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR200036-RT1488

Manual Integrations
 APPROVED

mohammad
 10/20/2019 8:16:44 PM

Quant Time: Oct 18 08:32:48 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	194859	20.00	ng	-0.02
21) Naphthalene-d8	8.03	136	690496	20.00	ng	-0.02
39) Acenaphthene-d10	9.80	164	366729	20.00	ng	-0.02
64) Phenanthrene-d10	11.29	188	664172	20.00	ng	-0.02
76) Chrysene-d12	13.97	240	547077	20.00	ng	-0.01
87) Perylene-d12	15.45	264	639041	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	1237668	102.10	ng	-0.01
7) Phenol-d6	6.39	99	1571307	107.57	ng	-0.02
23) Nitrobenzene-d5	7.31	82	879307	77.78	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	491663	125.81	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	1943154	85.73	ng	-0.02
79) Terphenyl-d14	12.90	244	2154662	79.74	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.51	163	258173	10.132	ng	Qvalue 100
71) Phenanthrene	11.32	178	95479	2.862	ng	99
75) Fluoranthene	12.52	202	258470	6.899	ng	99
78) Pyrene	12.75	202	220593	5.846	ng	99
81) Benzo(a)anthracene	13.96	228	83429	2.499	ng	95
83) Chrysene	14.00	228	86563	2.642	ng	98
88) Benzo(b)fluoranthene	15.02	252	105990m	2.924	ng	

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

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