

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012519\
 Data File : BM018654.D
 Acq On : 25 Jan 2019 16:07
 Operator : JU/SJ
 Sample : K1157-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-01-011519

Manual Integrations
 APPROVED

mohammad
 1/29/2019 12:27:03 AM

Quant Time: Jan 25 16:49:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 18 22:01:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	242933	20.00	ng	-0.01
21) Naphthalene-d8	10.52	136	938445	20.00	ng	0.00
39) Acenaphthene-d10	14.38	164	525004	20.00	ng	0.00
64) Phenanthrene-d10	17.13	188	1217209	20.00	ng	-0.01
76) Chrysene-d12	21.32	240	1328017	20.00	ng	-0.01
87) Perylene-d12	23.59	264	1332611	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	1830695	139.15	ng	0.00
7) Phenol-d6	6.91	99	2404641	143.90	ng	0.00
23) Nitrobenzene-d5	8.90	82	1465077	90.50	ng	0.00
42) 2,4,6-Tribromophenol	15.87	330	1027161	153.66	ng	-0.01
45) 2-Fluorobiphenyl	13.00	172	3276707	81.44	ng	0.00
79) Terphenyl-d14	19.76	244	4899199	77.77	ng	0.00
Target Compounds						
10) Phenol	6.93	94	41050	2.417	ng	92
50) Dimethylphthalate	13.84	163	304190	7.055	ng	99
75) Fluoranthene	19.19	202	223432	2.993	ng	98
78) Pyrene	19.56	202	240137	3.051	ng	100
88) Benzo(b)fluoranthene	22.91	252	164084m	2.169	ng	

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

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