

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM051019\
 Data File : BM020257.D
 Acq On : 11 May 2019 02:06
 Operator : JU/SJ
 Sample : K2801-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RT-3456

Quant Time: May 11 06:03:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 03:58:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	111333	20.00	ng	-0.02
21) Naphthalene-d8	10.57	136	427978	20.00	ng	-0.02
39) Acenaphthene-d10	14.42	164	277395	20.00	ng	-0.02
64) Phenanthrene-d10	17.17	188	664858	20.00	ng	-0.01
76) Chrysene-d12	21.35	240	517500	20.00	ng	-0.02
87) Perylene-d12	23.64	264	506474	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	708527	104.03	ng	-0.01
7) Phenol-d6	6.95	99	1000129	107.48	ng	0.00
23) Nitrobenzene-d5	8.94	82	689901	63.64	ng	-0.02
42) 2,4,6-Tribromophenol	15.92	330	477965	111.01	ng	0.00
45) 2-Fluorobiphenyl	13.04	172	1347577	59.77	ng	-0.02
79) Terphenyl-d14	19.79	244	2132375	71.87	ng	-0.02
Target Compounds						
50) Dimethylphthalate	13.87	163	185249	7.857	ng	98
75) Fluoranthene	19.23	202	120229	2.589	ng	98
78) Pyrene	19.59	202	113858	3.277	ng	99
83) Chrysene	21.39	228	84341	2.396	ng	100
88) Benzo(b)fluoranthene	22.95	252	70122	2.097	ng	# 88

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

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