

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030119\
 Data File : BM018995.D
 Acq On : 01 Mar 2019 20:08
 Operator : JU/SJ
 Sample : K1665-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 200031

Manual Integrations
 APPROVED

Sohil
 3/4/2019 4:53:01 PM

Quant Time: Mar 02 02:42:25 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 16:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	287434	20.00	ng	-0.04
21) Naphthalene-d8	10.47	136	1149008	20.00	ng	-0.04
39) Acenaphthene-d10	14.33	164	659589	20.00	ng	-0.03
64) Phenanthrene-d10	17.09	188	1420118	20.00	ng	-0.03
76) Chrysene-d12	21.28	240	1182068	20.00	ng	-0.03
87) Perylene-d12	23.53	264	1120749	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	2578483	146.98	ng	-0.02
7) Phenol-d6	6.86	99	3709502	150.10	ng	-0.02
23) Nitrobenzene-d5	8.85	82	2304412	92.73	ng	-0.04
42) 2,4,6-Tribromophenol	15.83	330	1425341	149.91	ng	-0.03
45) 2-Fluorobiphenyl	12.95	172	4575200	92.08	ng	-0.03
79) Terphenyl-d14	19.73	244	6013089	104.94	ng	-0.02
Target Compounds						
50) Dimethylphthalate	13.80	163	311867	5.673	ng	99
75) Fluoranthene	19.15	202	324125	3.677	ng	98
78) Pyrene	19.52	202	290119	4.226	ng	99
81) Benzo(a)anthracene	21.27	228	187794	2.725	ng	98
83) Chrysene	21.32	228	152816	2.314	ng	97
88) Benzo(b)fluoranthene	22.85	252	175984m	2.745	ng	

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

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