

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021819\
 Data File : BM018842.D
 Acq On : 18 Feb 2019 17:23
 Operator : JU/SJ
 Sample : K1518-17
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-MH-C02

Manual Integrations
 APPROVED

Sohil
 2/19/2019 10:03:10 AM

Quant Time: Feb 19 01:34:01 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	7.70	152	306738	20.00	ng	-0.01
21) Naphthalene-d8	10.49	136	1243280	20.00	ng	-0.01
39) Acenaphthene-d10	14.36	164	651299	20.00	ng	0.00
64) Phenanthrene-d10	17.10	188	1398728	20.00	ng	-0.01
76) Chrysene-d12	21.30	240	1088510	20.00	ng	0.00
87) Perylene-d12	23.56	264	1013724	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	2339765	124.98	ng	0.00
7) Phenol-d6	6.88	99	3262123	123.69	ng	0.00
23) Nitrobenzene-d5	8.87	82	2060102	76.62	ng	-0.01
42) 2,4,6-Tribromophenol	15.85	330	1081989	115.25	ng	-0.01
45) 2-Fluorobiphenyl	12.97	172	3978056	81.08	ng	-0.01
79) Terphenyl-d14	19.74	244	5043481	95.58	ng	0.00
Target Compounds						
10) Phenol	6.90	94	100371	3.617	ng	83
50) Dimethylphthalate	13.82	163	260135	4.793	ng	100
78) Pyrene	19.53	202	140635	2.225	ng	99
88) Benzo(b)fluoranthene	22.87	252	122641m	2.115	ng	

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

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