

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031519\
 Data File : BM019189.D
 Acq On : 15 Mar 2019 16:50
 Operator : JU/SJ
 Sample : K1908-03 10X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 101-102

Manual Integrations
 APPROVED

Sohil
 3/18/2019 3:04:33 PM

Quant Time: Mar 16 02:30:19 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 11 17:30:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	199772	20.00	ng	-0.01
21) Naphthalene-d8	10.44	136	934990	20.00	ng	-0.01
39) Acenaphthene-d10	14.31	164	562784	20.00	ng	-0.01
64) Phenanthrene-d10	17.06	188	1243003	20.00	ng	-0.01
76) Chrysene-d12	21.26	240	1159041	20.00	ng	-0.01
87) Perylene-d12	23.50	264	1032755	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	147719	11.98	ng	0.00
7) Phenol-d6	6.83	99	220314	12.21	ng	-0.01
23) Nitrobenzene-d5	8.83	82	127746	6.46	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	81645	9.83	ng	0.00
45) 2-Fluorobiphenyl	12.92	172	293239	6.78	ng	-0.01
79) Terphenyl-d14	19.70	244	427835	7.33	ng	-0.01
Target Compounds						
71) Phenanthrene	17.10	178	482592	6.624	ng	100
75) Fluoranthene	19.13	202	811135	9.702	ng	99
78) Pyrene	19.49	202	747422	10.192	ng	100
81) Benzo(a)anthracene	21.24	228	339699	4.672	ng	98
83) Chrysene	21.30	228	301019	4.279	ng	97
88) Benzo(b)fluoranthene	22.83	252	361866m	5.407	ng	
90) Benzo(a)pyrene	23.40	252	245561	4.083	ng	94
92) Benzo(g,h,i)perylene	26.45	276	137267	2.599	ng	97

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

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