

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080619\
 Data File : BM021865.D
 Acq On : 06 Aug 2019 19:17
 Operator : HP/JU
 Sample : K3863-18MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A52H9MS

Quant Time: Aug 06 19:52:22 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM080319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 06 01:29:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	702	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	3132	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.28	164	1850	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	3911	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	3858	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	4497	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	1150	0.22	ng/ul	0.01
14) Fluoranthene-d10	19.08	212	3128	0.24	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.35	153	1713	0.243	ng/ul	97
11) Pentachlorophenol	16.72	266	75	0.104	ng/ul	90
12) Phenanthrene	17.09	178	492	0.043	ng/ul#	75
15) Fluoranthene	19.11	202	1411	0.093	ng/ul	84
17) Pyrene	19.47	202	5187	0.327	ng/ul	96
18) Benzo(a)anthracene	21.23	228	575	0.042	ng/ul#	71
19) Chrysene	21.28	228	675	0.038	ng/ul#	82

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

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