

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080619\
 Data File : BM021866.D
 Acq On : 06 Aug 2019 19:54
 Operator : HP/JU
 Sample : K3863-19MSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A52H9MSD

Quant Time: Aug 06 20:27:24 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	7.64	152	714	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	3119	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.28	164	1859	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	3931	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	3850	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	4465	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	1078	0.21	ng/ul	0.02
14) Fluoranthene-d10	19.08	212	2945	0.22	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.35	153	1638	0.231	ng/ul	97
11) Pentachlorophenol	16.72	266	66	0.091	ng/ul	91
12) Phenanthrene	17.09	178	481	0.042	ng/ul#	70
15) Fluoranthene	19.11	202	1376	0.090	ng/ul	86
17) Pyrene	19.47	202	4906	0.310	ng/ul#	94
18) Benzo(a)anthracene	21.23	228	537	0.039	ng/ul#	75
19) Chrysene	21.29	228	697	0.039	ng/ul#	80

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

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