

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091619\
 Data File : BM022693.D
 Acq On : 16 Sep 2019 16:23
 Operator : HP/JU
 Sample : K4682-13MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0BL7MS

Quant Time: Sep 17 01:34:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM091019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 10 15:06:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	9327	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	37232	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	22738	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	48445	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	42741	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	43150	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	17425	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.23	212	47934	0.36	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.54	153	21603	0.245	ng/ul	99
11) Pentachlorophenol	16.86	266	2955	0.301	ng/ul	100
17) Pyrene	19.62	202	57769	0.298	ng/ul	98

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

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