

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM111616\
 Data File : BM007917.D
 Acq On : 16 Nov 2016 12:45
 Operator : UM/SJ
 Sample : H5622-10MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4WK9MS

Manual Integrations
 APPROVED

umangi
 11/17/2016 5:55:14 PM

Quant Time: Nov 17 06:36:49 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM102516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 16 16:16:41 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	961	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	3138	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.37	164	1788	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.11	188	3577m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.31	240	3380	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	2875m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.13	152	1114	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.16	212	2750m	0.34	ng/ul	0.00
Target Compounds						Qvalue
8) Acenaphthene	14.42	153	1682	0.26	ng/ul	91
11) Pentachlorophenol	16.80	266	309	0.37	ng/ul	96
17) Pyrene	19.53	202	4523	0.39	ng/ul	97

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

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