

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112316\
 Data File : BM008079.D
 Acq On : 24 Nov 2016 18:03
 Operator : UM/SJ
 Sample : H5701-18
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4WF0

Manual Integrations
 APPROVED

Umangi
 11/28/2016 5:30:26 PM

Quant Time: Nov 25 03:51:30 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 25 01:27:01 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	727	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.49	136	2546	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.35	164	1592	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	3066m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.29	240	2939	0.40	ng/ul	0.00
20) Perylene-d12	23.52	264	2600m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	1403	0.39	ng/ul	-0.01
14) Fluoranthene-d10	19.13	212	3168	0.37	ng/ul	0.00
Target Compounds						
15) Fluoranthene	19.16	202	1796	0.15	ng/ul	96
17) Pyrene	19.52	202	2119	0.21	ng/ul#	95
18) Benzo(a)anthracene	21.27	228	693	0.08	ng/ul#	79
19) Chrysene	21.33	228	798	0.07	ng/ul#	82
23) Benzo(a)pyrene	23.42	252	707m	0.07	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.79	276	487m	0.04	ng/ul	
26) Benzo(g,h,i)perylene	26.47	276	568m	0.05	ng/ul	

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

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