

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092216\
 Data File : BM007512.D
 Acq On : 22 Sep 2016 15:40
 Operator : UM/SJ
 Sample : H4830-04MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4VH2MSD

Manual Integrations
 APPROVED

umangi
 9/23/2016 6:52:32 PM

Quant Time: Sep 22 17:17:33 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM092116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 22 01:26:49 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	1431	0.40	ng/ul	0.00
2) Naphthalene-d8	10.57	136	5793	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.42	164	4159	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.16	188	8637m	0.40	ng/ul	0.00
16) Chrysene-d12	21.35	240	9149	0.40	ng/ul	0.00
20) Perylene-d12	23.61	264	6884m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.17	152	3071	0.45	ng/ul	0.00
14) Fluoranthene-d10	19.19	212	9045	0.45	ng/ul	0.00
Target Compounds						
8) Acenaphthene	14.47	153	5353	0.39	ng/ul	88
11) Pentachlorophenol	16.83	266	1123	0.48	ng/ul	99
15) Fluoranthene	19.22	202	726	0.02	ng/ul	74
17) Pyrene	19.59	202	15156	0.50	ng/ul	98

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

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