

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092215\
 Data File : BM002702.D
 Acq On : 22 Sep 2015 20:00
 Operator : UM/IZ
 Sample : G3741-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHAF4

Manual Integrations
 APPROVED

MMDadoda
 9/23/2015 4:37:27 PM

Quant Time: Sep 23 09:02:04 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM90315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 23 03:40:59 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.78	152	9185	0.40	ng/ul	0.00
4) Naphthalene-d8	11.64	136	35291	0.40	ng/ul	0.00
8) Acenaphthene-d10	15.36	164	25346	0.40	ng/ul	0.00
12) Phenanthrene-d10	18.06	188	31383	0.40	ng/ul	0.00
18) Chrysene-d12	22.16	240	27187	0.40	ng/ul	0.00
22) Perylene-d12	24.77	264	27142	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
6) 2-Methylnaphthalene-d10	13.16	152	16734m	0.35	ng/ul	0.00
16) Fluoranthene-d10	20.03	212	21500	0.30	ng/ul	0.00
Target Compounds						
17) Fluoranthene	20.06	202	15360	0.16	ng/ul	81
19) Pyrene	20.42	202	43387	0.42	ng/ul#	90
20) Benzo(a)anthracene	22.13	228	5708m	0.06	ng/ul	
21) Chrysene	22.19	228	7353m	0.08	ng/ul	
23) Benzo(b)fluoranthene	23.96	252	8293m	0.09	ng/ul	
24) Benzo(k)fluoranthene	24.00	252	3262m	0.04	ng/ul	
25) Benzo(a)pyrene	24.66	252	5152	0.06	ng/ul#	53
26) Indeno(1,2,3-cd)pyrene	27.54	276	3301	0.03	ng/ul	95
28) Benzo(g,h,i)perylene	28.39	276	3201	0.04	ng/ul#	70

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

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