

Data Path : U:\HPCHEM1\BNA M\Data\BM092415\
 Data File : BM002716.D
 Acq On : 24 Sep 2015 17:39
 Operator : UM/IZ
 Sample : G3741-06MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHAF7MS

Manual Integrations
 APPROVED

MMDadoda
 9/25/2015 2:19:08 PM

Quant Time: Oct 02 11:49:46 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM092415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 25 02:23:28 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.74	152	13183	0.40	ng/ul	0.00
4) Naphthalene-d8	11.60	136	51884	0.40	ng/ul	0.00
8) Acenaphthene-d10	15.33	164	29678m	0.40	ng/ul	0.00
12) Phenanthrene-d10	18.03	188	61694	0.40	ng/ul	0.00
18) Chrysene-d12	22.12	240	53945	0.40	ng/ul	0.00
22) Perylene-d12	24.72	264	50100	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	13.13	152	21972m	0.29	ng/ul	0.00
16) Fluoranthene-d10	20.00	212	50532	0.33	ng/ul	-0.01
Target Compounds				Ovalue		
5) Naphthalene	11.64	128	209880	1.49	ng/ul#	61
7) 2-Methylnaphthalene	13.20	142	335990	3.71	ng/ul	98
10) Acenaphthene	15.38	153	103832	0.93	ng/ul	98
11) Fluorene	16.36	166	314654	2.47	ng/ul	86
13) Pentachlorophenol	17.69	266	5614	0.47	ng/ul	97
14) Phenanthrene	18.08	178	152020	0.76	ng/ul#	94
17) Fluoranthene	20.04	202	15868	0.07	ng/ul	69
19) Pyrene	20.39	202	84080	0.40	ng/ul	96
20) Benzo(a)anthracene	22.10	228	4795	0.03	ng/ul#	70
21) Chrysene	22.16	228	4783	0.03	ng/ul#	74
26) Indeno(1,2,3-cd)pyrene	27.45	276	4022	0.02	ng/ul	94

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

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