

Data Path : Z:\HPCHEM1\BNA_M\Data\BM092815\
 Data File : BM002731.D
 Acq On : 28 Sep 2015 12:55
 Operator : UM/IZ
 Sample : G3798-06
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9G60

Manual Integrations
 APPROVED

MMDadoda
 9/29/2015 3:42:12 PM

Quant Time: Oct 13 05:39:28 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 : Tue Sep 29 00:32:23 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.74	152	9900	0.40	ng/ul	0.00
4) Naphthalene-d8	11.60	136	38476	0.40	ng/ul	0.01
8) Acenaphthene-d10	15.33	164	20601	0.40	ng/ul	0.00
12) Phenanthrene-d10	18.03	188	44063	0.40	ng/ul	0.00
18) Chrysene-d12	22.13	240	44065	0.40	ng/ul	0.00
22) Perylene-d12	24.72	264	39845	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.80	96	46832	4.58	ng/uL	0.00
6) 2-Methylnaphthalene-d10	13.13	152	14487	0.26	ng/ul	0.00
16) Fluoranthene-d10	20.00	212	28330	0.26	ng/ul	0.00
Target Compounds						
					Qvalue	
7) 2-Methylnaphthalene	13.20	142	532	0.01	ng/ul#	71
14) Phenanthrene	18.08	178	4887	0.03	ng/ul#	91
17) Fluoranthene	20.03	202	6949	0.04	ng/ul	91
19) Pyrene	20.39	202	6755	0.04	ng/ul#	88
20) Benzo(a)anthracene	22.10	228	2450	0.02	ng/ul#	76
21) Chrysene	22.16	228	3400	0.02	ng/ul#	80
23) Benzo(b)fluoranthene	23.92	252	4477m	0.03	ng/ul	
28) Benzo(g,h,i)perylene	28.31	276	2204	0.02	ng/ul#	57

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

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