

Data Path : Z:\HPCHEM1\BNA E\DATA\BE111115\
 Data File : BE091140.D
 Acq On : 12 Nov 2015 12:05
 Operator : UM/NP
 Sample : G4322-13
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A4HS4

Manual Integrations
 APPROVED

Mmdadoda
 11/16/2015 12:42:31 PM

Quant Time: Nov 13 16:45:56 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM02.2-EPA-SIM-BE102715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 13 13:56:00 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	3653	0.40	ng/ul	0.00
4) Naphthalene-d8	9.57	136	20535	0.40	ng/ul	0.00
8) Acenaphthene-d10	13.40	164	12280	0.40	ng/ul	0.00
12) Phenanthrene-d10	16.12	188	29715m	0.40	ng/ul	-0.01
18) Chrysene-d12	20.28	240	40510	0.40	ng/ul	0.00
22) Perylene-d12	22.19	264	44905	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	2.93	96	9431	3.59	ng/uL	-0.02
6) 2-Methylnaphthalene-d10	11.07	152	3588	0.13	ng/ul	0.00
16) Fluoranthene-d10	18.12	212	10306	0.12	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	9.61	128	1966	0.04	ng/ul#	85
9) Acenaphthylene	13.14	152	9183	0.15	ng/ul	98
11) Fluorene	14.45	166	2272	0.04	ng/ul#	92
14) Phenanthrene	16.16	178	230137	0.67	ng/ul	97
15) Anthracene	16.24	178	52532m	0.17	ng/ul	
17) Fluoranthene	18.15	202	528377m	1.32	ng/ul	
19) Pyrene	18.53	202	528780	1.15	ng/ul#	80
20) Benzo(a)anthracene	20.26	228	54990m	0.53	ng/ul	
21) Chrysene	20.32	228	72569	0.61	ng/ul	97
23) Benzo(b)fluoranthene	21.62	252	56780	0.41	ng/ul	97
24) Benzo(k)fluoranthene	21.65	252	59760	0.42	ng/ul	95
25) Benzo(a)pyrene	22.09	252	60854	0.47	ng/ul	94
26) Indeno(1,2,3-cd)pyrene	23.83	276	164877m	0.30	ng/ul	
27) Dibenzo(a,h)anthracene	23.79	278	10296	0.08	ng/ul#	65
28) Benzo(a,h,i)perylene	24.40	276	163888	0.29	ng/ul	97

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

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