

Data Path : Z:\HPCHEM1\BNA E\DATA\BE111115\
 Data File : BE091132.D
 Acq On : 12 Nov 2015 7:13
 Operator : UM/NP
 Sample : G4322-04
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A4HR5

Manual Integrations
 APPROVED

Mmdadoda
 11/16/2015 12:42:18 PM

Quant Time: Nov 13 18:05:36 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	6.90	152	3668	0.40	ng/ul	0.00
4) Naphthalene-d8	9.57	136	20214	0.40	ng/ul	0.00
8) Acenaphthene-d10	13.40	164	11435	0.40	ng/ul	0.00
12) Phenanthrene-d10	16.12	188	29224m	0.40	ng/ul	-0.01
18) Chrysene-d12	20.28	240	38741	0.40	ng/ul	0.00
22) Perylene-d12	22.19	264	34219m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	2.92	96	11334	4.30	ng/uL	-0.03
6) 2-Methylnaphthalene-d10	11.07	152	3065	0.11	ng/ul	0.00
16) Fluoranthene-d10	18.12	212	6959	0.08	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	9.61	128	3524	0.07	ng/ul#	92
9) Acenaphthylene	13.15	152	4573	0.08	ng/ul	97
10) Acenaphthene	13.46	153	1309	0.03	ng/ul	88
11) Fluorene	14.45	166	1695	0.03	ng/ul#	90
14) Phenanthrene	16.16	178	194534	0.58	ng/ul	98
15) Anthracene	16.25	178	39005m	0.13	ng/ul	
17) Fluoranthene	18.15	202	440317	1.11	ng/ul	88
19) Pyrene	18.53	202	416714	0.95	ng/ul#	80
20) Benzo(a)anthracene	20.26	228	40736	0.41	ng/ul	96
21) Chrysene	20.31	228	56590	0.49	ng/ul	97
23) Benzo(b)fluoranthene	21.62	252	47735	0.45	ng/ul	94
24) Benzo(k)fluoranthene	21.65	252	49718	0.45	ng/ul	96
25) Benzo(a)pyrene	22.09	252	50733	0.52	ng/ul	92
26) Indeno(1,2,3-cd)pyrene	23.83	276	135596m	0.33	ng/ul	
27) Dibenzo(a,h)anthracene	23.78	278	7560	0.08	ng/ul#	73
28) Benzo(a,h,i)perylene	24.39	276	134067	0.31	ng/ul	98

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

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