

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE111515\
 Data File : BE091167.D
 Acq On : 15 Nov 2015 11:39
 Operator : UM/NP
 Sample : G4322-06DL 5X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A4HR7DL

Manual Integrations
 APPROVED

mmdadoda
 11/16/2015 3:27:04 PM

Quant Time: Nov 16 06:33:18 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 : Mon Nov 16 04:54:02 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	7208	0.40	ng/ul	-0.03
4) Naphthalene-d8	9.58	136	35242	0.40	ng/ul	0.00
8) Acenaphthene-d10	13.39	164	20577	0.40	ng/ul	0.00
12) Phenanthrene-d10	16.10	188	52022	0.40	ng/ul	0.00
18) Chrysene-d12	20.23	240	87258	0.40	ng/ul	0.00
22) Perylene-d12	22.12	264	83034	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.16	96	2419m	0.47	ng/uL	0.07
6) 2-Methylnaphthalene-d10	11.11	152	451	0.01	ng/ul	0.02
16) Fluoranthene-d10	18.08	212	1173	0.01	ng/ul	0.00
Target Compounds				Qvalue		
5) Naphthalene	9.62	128	17414	0.20	ng/ul	98
7) 2-Methylnaphthalene	11.18	142	1754	0.03	ng/ul	96
14) Phenanthrene	16.14	178	134980	0.22	ng/ul	96
15) Anthracene	16.22	178	27036m	0.05	ng/ul	
17) Fluoranthene	18.11	202	180148	0.26	ng/ul	87
19) Pyrene	18.49	202	152972	0.15	ng/ul#	78
20) Benzo(a)anthracene	20.21	228	21419	0.10	ng/ul#	94
21) Chrysene	20.27	228	26610	0.10	ng/ul#	96
23) Benzo(b)fluoranthene	21.57	252	18334	0.07	ng/ul	92
24) Benzo(k)fluoranthene	21.59	252	21410	0.08	ng/ul#	88
25) Benzo(a)pyrene	22.02	252	20372	0.09	ng/ul#	93
26) Indeno(1,2,3-cd)pyrene	23.73	276	54128	0.05	ng/ul#	91
28) Benzo(g,h,i)perylene	24.28	276	51548	0.05	ng/ul	96

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

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