

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112315\
 Data File : BE091305.D
 Acq On : 23 Nov 2015 18:22
 Operator : UM/NP
 Sample : G4322-04DL 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A4HR5DL

Quant Time: Nov 24 03:02:07 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 : Tue Nov 24 02:51:08 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	1942	0.40	ng/ul	-0.02
4) Naphthalene-d8	9.57	136	10330	0.40	ng/ul	0.00
8) Acenaphthene-d10	13.37	164	5624	0.40	ng/ul	0.00
12) Phenanthrene-d10	16.09	188	16532	0.40	ng/ul	0.00
18) Chrysene-d12	20.23	240	21358	0.40	ng/ul	0.00
22) Perylene-d12	22.11	264	20220	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	2.97	96	2584	1.85	ng/uL	0.02
6) 2-Methylnaphthalene-d10	11.09	152	444	0.03	ng/ul	0.02
16) Fluoranthene-d10	18.07	212	1234	0.03	ng/ul	0.00
Target Compounds						Ovalue
14) Phenanthrene	16.12	178	28241	0.15	ng/ul	96
17) Fluoranthene	18.10	202	70047	0.31	ng/ul	90
19) Pyrene	18.47	202	68248	0.28	ng/ul#	80
20) Benzo(a)anthracene	20.20	228	5286	0.10	ng/ul	97
21) Chrysene	20.26	228	9461	0.15	ng/ul	97
23) Benzo(b)fluoranthene	21.55	252	7010	0.11	ng/ul	97
24) Benzo(k)fluoranthene	21.58	252	7410	0.11	ng/ul	97
25) Benzo(a)pyrene	22.01	252	7731	0.13	ng/ul	98
26) Indeno(1,2,3-cd)pyrene	23.71	276	17653	0.07	ng/ul#	90
28) Benzo(g,h,i)perylene	24.26	276	18188	0.07	ng/ul	98

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

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