

Data Path : Z:\HPCHEM1\BNA E\DATA\BE111615\
 Data File : BE091182.D
 Acq On : 16 Nov 2015 13:54
 Operator : UM/NP
 Sample : G4309-09
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Nov 17 10:34:52 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	6.88	152	4035	0.40	ng/ul	-0.05
4) Naphthalene-d8	9.56	136	20618	0.40	ng/ul	-0.03
8) Acenaphthene-d10	13.37	164	12496	0.40	ng/ul	-0.01
12) Phenanthrene-d10	16.17	188	1542370	0.40	ng/ul	0.07
18) Chrysene-d12	20.23	240	54748	0.40	ng/ul	0.00
22) Perylene-d12	22.12	264	66564	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.02	96	21404	7.37	ng/uL	-0.07
6) 2-Methylnaphthalene-d10	11.07	152	7054	0.25	ng/ul	-0.02
16) Fluoranthene-d10	18.08	212	21916	0.01	ng/ul	0.00
Target Compounds				Ovalue		
3) 1,4-Dioxane	3.18	88	179	0.05	ng/ul#	42
19) Pyrene	18.48	202	103148	0.17	ng/ul#	78
20) Benzo(a)anthracene	20.26	228	14532	0.10	ng/ul	95
21) Chrysene	20.26	228	14797	0.09	ng/ul	98
23) Benzo(b)fluoranthene	21.56	252	16756	0.08	ng/ul	94
24) Benzo(k)fluoranthene	21.59	252	9120	0.04	ng/ul#	89
25) Benzo(a)pyrene	22.02	252	11978	0.06	ng/ul#	93
26) Indeno(1,2,3-cd)pyrene	23.72	276	34324	0.04	ng/ul#	90
28) Benzo(g,h,i)perylene	24.27	276	40344	0.05	ng/ul	93

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

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