

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE111715\
 Data File : BE091216.D
 Acq On : 17 Nov 2015 22:22
 Operator : UM/NP
 Sample : G4273-05DL 25X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AP0DL

Quant Time: Nov 18 04:27:06 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 : Wed Nov 18 04:23:52 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	7	0.40	ng/ul	0.00
4) Naphthalene-d8	9.57	136	6	0.40	ng/ul	0.00
8) Acenaphthene-d10	13.42	164	16	0.40	ng/ul	0.04
12) Phenanthrene-d10	16.14	188	26	0.40	ng/ul	0.05
18) Chrysene-d12	20.28	240	31	0.40	ng/ul	0.05
22) Perylene-d12	22.13	264	12	0.40	ng/ul	0.02
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.02	96	24	4.77	ng/uL	0.00
6) 2-Methylnaphthalene-d10	11.18	152	5	0.61	ng/ul	0.12
16) Fluoranthene-d10	18.13	212	3	0.04	ng/ul	0.06
Target Compounds				Qvalue		
3) 1,4-Dioxane	3.06	88	18	3.16	ng/ul#	81
7) 2-Methylnaphthalene	11.12	142	7	0.73	ng/ul	82
9) Acenaphthylene	13.16	152	16	0.20	ng/ul#	1
10) Acenaphthene	13.50	153	10	0.18	ng/ul#	62
11) Fluorene	14.44	166	3	0.04	ng/ul#	68
13) Pentachlorophenol	15.91	266	1	0.40	ng/ul#	16
14) Phenanthrene	16.20	178	389	1.30	ng/ul#	1
15) Anthracene	16.20	178	389	1.42	ng/ul#	1
17) Fluoranthene	18.15	202	98	0.28	ng/ul#	1
19) Pyrene	18.51	202	111	0.32	ng/ul#	18
20) Benzo(a)anthracene	20.24	228	20	0.25	ng/ul#	1
21) Chrysene	20.31	228	11	0.12	ng/ul#	1
23) Benzo(b)fluoranthene	21.58	252	34	0.91	ng/ul#	1
24) Benzo(k)fluoranthene	21.62	252	27	0.70	ng/ul#	1
25) Benzo(a)pyrene	22.04	252	21	0.61	ng/ul#	1
26) Indeno(1,2,3-cd)pyrene	23.75	276	107	0.74	ng/ul#	28
27) Dibenzo(a,h)anthracene	23.70	278	22	0.64	ng/ul#	1
28) Benzo(g,h,i)perylene	24.28	276	58	0.38	ng/ul#	1

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

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