

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM091115\
 Data File : BM002680.D
 Acq On : 11 Sep 2015 10:28
 Operator : UM/IZ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Sep 12 01:29:30 2015

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM90315.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Sep 04 04:35:12 2015

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.78	152	8884	0.40	ng/ul	0.00
4) Naphthalene-d8	11.64	136	37156	0.40	ng/ul	0.00
8) Acenaphthene-d10	15.36	164	16976	0.40	ng/ul	0.00
12) Phenanthrene-d10	18.06	188	31439	0.40	ng/ul	-0.02
18) Chrysene-d12	22.16	240	25410	0.40	ng/ul	0.00
22) Perylene-d12	24.77	264	24196	0.40	ng/ul	-0.01

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.82	96	7261	0.81	ng/uL	0.00
6) 2-Methylnaphthalene-d10	13.17	152	18887	0.38	ng/ul	0.00
16) Fluoranthene-d10	20.03	212	26832	0.37	ng/ul	-0.01

Target Compounds

						Qvalue
3) 1,4-Dioxane	3.87	88	8392	0.81	ng/ul	98
5) Naphthalene	11.69	128	36364	0.38	ng/ul	98
7) 2-Methylnaphthalene	13.23	142	24473	0.38	ng/ul	99
9) Acenaphthylene	15.09	152	33627	0.37	ng/ul	95
10) Acenaphthene	15.42	153	24108	0.38	ng/ul	91
11) Fluorene	16.39	166	25479	0.38	ng/ul	86
13) Pentachlorophenol	17.72	266	951	0.31	ng/ul	97
14) Phenanthrene	18.11	178	35683	0.38	ng/ul	97
15) Anthracene	18.20	178	33858	0.36	ng/ul	97
17) Fluoranthene	20.06	202	36295	0.37	ng/ul	96
19) Pyrene	20.42	202	37181	0.39	ng/ul	97
20) Benzo(a)anthracene	22.14	228	31381	0.37	ng/ul	95
21) Chrysene	22.19	228	30636	0.38	ng/ul	99
23) Benzo(b)fluoranthene	23.96	252	33140	0.39	ng/ul	97
24) Benzo(k)fluoranthene	24.01	252	29409	0.37	ng/ul	95
25) Benzo(a)pyrene	24.66	252	29576	0.37	ng/ul	96
26) Indeno(1,2,3-cd)pyrene	27.54	276	32202	0.34	ng/ul	98
27) Dibenzo(a,h)anthracene	27.55	278	24835	0.32	ng/ul	94
28) Benzo(g,h,i)perylene	28.40	276	28945	0.36	ng/ul	96

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

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