

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042715\
 Data File : BM001114.D
 Acq On : 27 Apr 2015 11:53
 Operator : TP/IZ
 Sample : SSTD0.445
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.445

Quant Time: Apr 27 14:12:03 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM042715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 27 13:50:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	974	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	3344	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	1707	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	3843	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	3523	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	3435	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.93	152	1714	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	3464	0.41	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.37	128	3462	0.40	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	2255	0.39	ng/ul	100
7) Acenaphthylene	13.91	152	3113	0.41	ng/ul	100
8) Acenaphthene	14.27	153	2477	0.41	ng/ul	100
9) Fluorene	15.26	166	2766	0.40	ng/ul	100
11) Pentachlorophenol	16.63	266	248	0.29	ng/ul	100
12) Phenanthrene	17.00	178	4375	0.39	ng/ul	100
13) Anthracene	17.09	178	4053	0.40	ng/ul	100
15) Fluoranthene	19.02	202	5084	0.41	ng/ul	100
17) Pyrene	19.38	202	5269	0.39	ng/ul	100
18) Benzo(a)anthracene	21.14	228	4544	0.41	ng/ul	100
19) Chrysene	21.19	228	5029	0.41	ng/ul	100
21) Benzo(b)fluoranthene	22.67	252	4921	0.39	ng/ul	100
22) Benzo(k)fluoranthene	22.71	252	5079	0.40	ng/ul	100
23) Benzo(a)pyrene	23.22	252	4849	0.41	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.46	276	5561	0.42	ng/ul	100
25) Dibenzo(a,h)anthracene	25.49	278	4446	0.41	ng/ul	100
26) Benzo(g,h,i)perylene	26.12	276	4853	0.41	ng/ul	100

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

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