

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031215\
 Data File : BM000701.D
 Acq On : 12 Mar 2015 16:17
 Operator : TP/IZ
 Sample : SSTD1.675
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.675

Quant Time: Mar 13 09:02:33 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM021815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 13 08:51:18 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	3816	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	13595	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	1998	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	17372	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	17649	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	16321	0.40	ng/ul	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.02	152	28620	1.55	ng/ul	-0.01
14) Fluoranthene-d10	19.08	212	64212	1.68	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.48	128	57058	1.62	ng/ul	99
5) 2-Methylnaphthalene	12.10	142	39477	1.59	ng/ul	99
7) Acenaphthylene	14.02	152	56827	4.18	ng/ul	99
8) Acenaphthene	14.36	153	42332	4.05	ng/ul	99
9) Fluorene	15.35	166	52461	4.08	ng/ul	97
11) Pentachlorophenol	16.71	266	3982	0.89	ng/ul	98
12) Phenanthrene	17.08	178	81937	1.61	ng/ul	99
13) Anthracene	17.17	178	78658	1.65	ng/ul	99
15) Fluoranthene	19.11	202	98806	1.70	ng/ul	98
17) Pyrene	19.48	202	100009	1.26	ng/ul	99
18) Benzo(a)anthracene	21.23	228	89060	1.57	ng/ul	99
19) Chrysene	21.28	228	99033	1.60	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	99448	1.59	ng/ul	97
22) Benzo(k)fluoranthene	22.82	252	94076	1.53	ng/ul	97
23) Benzo(a)pyrene	23.34	252	90156	1.61	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.65	276	104960	1.94	ng/ul	100
25) Dibenzo(a,h)anthracene	25.66	278	84714	1.97	ng/ul	97
26) Benzo(g,h,i)perylene	26.31	276	91209	1.94	ng/ul	98

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

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