

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031915\
 Data File : BM000736.D
 Acq On : 19 Mar 2015 21:08
 Operator : TP/IZ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.402

Quant Time: Mar 20 11:34:50 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM031915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 19 15:39:49 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	3025	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	10204	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	5201	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	9764	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	8089	0.40	ng/ul	0.01
20) Perylene-d12	23.41	264	7423	0.40	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.00	152	5251	0.45	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	8679	0.45	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.46	128	11141	0.46	ng/ul	100
5) 2-Methylnaphthalene	12.08	142	7042	0.45	ng/ul	100
7) Acenaphthylene	14.00	152	8738	0.48	ng/ul	99
8) Acenaphthene	14.33	153	7027	0.48	ng/ul	98
9) Fluorene	15.34	166	7764	0.47	ng/ul	97
11) Pentachlorophenol	16.68	266	407	0.33	ng/ul	94
12) Phenanthrene	17.05	178	11816	0.45	ng/ul	100
13) Anthracene	17.15	178	10442	0.46	ng/ul	99
15) Fluoranthene	19.09	202	12801	0.44	ng/ul	99
17) Pyrene	19.45	202	12976	0.45	ng/ul	100
18) Benzo(a)anthracene	21.21	228	9994	0.44	ng/ul	91
19) Chrysene	21.26	228	12206	0.46	ng/ul#	90
21) Benzo(b)fluoranthene	22.75	252	11227	0.44	ng/ul	98
22) Benzo(k)fluoranthene	22.79	252	12262	0.47	ng/ul	99
23) Benzo(a)pyrene	23.31	252	10526	0.46	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.60	276	12144	0.45	ng/ul	100
25) Dibenzo(a,h)anthracene	25.62	278	9810	0.45	ng/ul	98
26) Benzo(g,h,i)perylene	26.26	276	10854	0.46	ng/ul	99

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

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