

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031915\
 Data File : BM000728.D
 Acq On : 19 Mar 2015 14:01
 Operator : TP/IZ
 Sample : SSTD0.898
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.898

Quant Time: Mar 19 15:13:44 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 14:57:23 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	2842	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	9715	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	5398	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	9675	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	8120	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	7256	0.40	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.00	152	9448	0.87	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	15871	0.88	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.46	128	19637	0.85	ng/ul	99
5) 2-Methylnaphthalene	12.08	142	12466	0.86	ng/ul	100
7) Acenaphthylene	14.00	152	15809	0.79	ng/ul	98
8) Acenaphthene	14.33	153	12475	0.76	ng/ul	99
9) Fluorene	15.34	166	14227	0.78	ng/ul	97
11) Pentachlorophenol	16.68	266	792	0.91	ng/ul	96
12) Phenanthrene	17.05	178	21712	0.85	ng/ul	99
13) Anthracene	17.15	178	19378	0.88	ng/ul	98
15) Fluoranthene	19.09	202	24247	0.90	ng/ul	99
17) Pyrene	19.45	202	24147	0.80	ng/ul	99
18) Benzo(a)anthracene	21.20	228	19129	0.87	ng/ul	100
19) Chrysene	21.25	228	22504	0.84	ng/ul	98
21) Benzo(b)fluoranthene	22.75	252	20844	0.81	ng/ul	98
22) Benzo(k)fluoranthene	22.79	252	20949	0.81	ng/ul	97
23) Benzo(a)pyrene	23.31	252	19259	0.85	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.60	276	22465	0.86	ng/ul	100
25) Dibenzo(a,h)anthracene	25.61	278	18182	0.86	ng/ul	94
26) Benzo(g,h,i)perylene	26.26	276	19888	0.85	ng/ul	98

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

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