

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031215\
 Data File : BM000698.D
 Acq On : 12 Mar 2015 14:30
 Operator : TP/IZ
 Sample : SSTD0.272
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.272

Quant Time: Mar 13 09:01:51 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	4366	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	15509	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	2425	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	19293	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	17381	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	15835	0.40	ng/ul	0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	4092	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	8576	0.20	ng/ul	0.01

Target Compounds				Qvalue		
3) Naphthalene	10.48	128	8083	0.20	ng/ul	97
5) 2-Methylnaphthalene	12.11	142	5464	0.19	ng/ul	99
7) Acenaphthylene	14.02	152	7471	0.45	ng/ul	99
8) Acenaphthene	14.36	153	5458	0.43	ng/ul	99
9) Fluorene	15.35	166	6959	0.45	ng/ul	97
11) Pentachlorophenol	16.71	266	460	0.09	ng/ul	93
12) Phenanthrene	17.08	178	11219	0.20	ng/ul	99
13) Anthracene	17.17	178	10083	0.19	ng/ul	98
15) Fluoranthene	19.11	202	12840	0.20	ng/ul	98
17) Pyrene	19.48	202	13363	0.17	ng/ul	100
18) Benzo(a)anthracene	21.23	228	10426	0.19	ng/ul	99
19) Chrysene	21.28	228	12786	0.21	ng/ul	98
21) Benzo(b)fluoranthene	22.78	252	11366	0.19	ng/ul	94
22) Benzo(k)fluoranthene	22.82	252	11815	0.20	ng/ul	94
23) Benzo(a)pyrene	23.34	252	10707	0.20	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.65	276	11083	0.21	ng/ul	99
25) Dibenzo(a,h)anthracene	25.66	278	8710	0.21	ng/ul	94
26) Benzo(g,h,i)perylene	26.32	276	9839	0.22	ng/ul	96

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

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