

Data Path : Z:\HPCHEM1\BNA_M\Data\BM071415\
 Data File : BM001917.D
 Acq On : 14 Jul 2015 16:43
 Operator : TP/UM
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.449

Quant Time: Jul 14 18:38:39 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM071415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 14 04:27:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	1374	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	5317	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	2814	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	6044	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	5174	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	4503	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.03	152	3368	0.41	ng/ul	-0.01
14) Fluoranthene-d10	19.09	212	6622	0.40	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.49	128	6506	0.40	ng/ul	100
5) 2-Methylnaphthalene	12.11	142	4333	0.40	ng/ul	99
7) Acenaphthylene	14.01	152	6627	0.40	ng/ul	99
8) Acenaphthene	14.37	153	4685	0.40	ng/ul	98
9) Fluorene	15.37	166	5446	0.40	ng/ul	96
11) Pentachlorophenol	16.70	266	588	0.37	ng/ul	95
12) Phenanthrene	17.09	178	8440	0.40	ng/ul	99
13) Anthracene	17.18	178	8104	0.40	ng/ul	100
15) Fluoranthene	19.12	202	9279	0.40	ng/ul	99
17) Pyrene	19.48	202	9486	0.40	ng/ul	99
18) Benzo(a)anthracene	21.23	228	8016	0.40	ng/ul	97
19) Chrysene	21.28	228	8091	0.40	ng/ul	99
21) Benzo(b)fluoranthene	22.81	252	7802	0.38	ng/ul	98
22) Benzo(k)fluoranthene	22.85	252	7186	0.42	ng/ul	99
23) Benzo(a)pyrene	23.38	252	6960	0.39	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.71	276	7051	0.37	ng/ul	99
25) Dibenzo(a,h)anthracene	25.74	278	5583	0.37	ng/ul	98
26) Benzo(g,h,i)perylene	26.40	276	5986	0.37	ng/ul	98

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

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