

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021615\
 Data File : BM000535.D
 Acq On : 16 Feb 2015 17:01
 Operator : TP/IZ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.414

Quant Time: Feb 17 02:26:18 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM021615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 16 14:03:19 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	3169	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	10656	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	6813	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	12465	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	10578	0.40	ng/ul	0.00
20) Perylene-d12	23.52	264	9043	0.40	ng/ul	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.09	152	5633	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	10868	0.39	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.54	128	11612	0.41	ng/ul	100
5) 2-Methylnaphthalene	12.16	142	7742	0.41	ng/ul	100
7) Acenaphthylene	14.07	152	10531	0.39	ng/ul	100
8) Acenaphthene	14.41	153	8356	0.40	ng/ul	98
9) Fluorene	15.40	166	9749	0.39	ng/ul	99
11) Pentachlorophenol	16.76	266	369	0.29	ng/ul	92
12) Phenanthrene	17.13	178	15580	0.40	ng/ul	99
13) Anthracene	17.23	178	13693	0.40	ng/ul	99
15) Fluoranthene	19.16	202	17608	0.40	ng/ul	89
17) Pyrene	19.53	202	17808	0.41	ng/ul	100
18) Benzo(a)anthracene	21.27	228	13349	0.40	ng/ul	94
19) Chrysene	21.32	228	15415	0.40	ng/ul	95
21) Benzo(b)fluoranthene	22.84	252	13739	0.40	ng/ul	99
22) Benzo(k)fluoranthene	22.89	252	14237	0.40	ng/ul	97
23) Benzo(a)pyrene	23.42	252	12631	0.39	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.75	276	13775	0.41	ng/ul	99
25) Dibenzo(a,h)anthracene	25.77	278	10778	0.41	ng/ul	98
26) Benzo(g,h,i)perylene	26.44	276	12166	0.40	ng/ul	99

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

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