

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021615\
 Data File : BM000529.D
 Acq On : 16 Feb 2015 11:30
 Operator : TP/IZ
 Sample : SSTD0.410
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Feb 16 13:17:28 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 13:07:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	3539	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	12202	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	7758	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	14977	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	13029	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	10866	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.09	152	6447	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	13036	0.40	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.54	128	13245	0.42	ng/ul	100
5) 2-Methylnaphthalene	12.16	142	8884	0.40	ng/ul	100
7) Acenaphthylene	14.07	152	12045	0.37	ng/ul	100
8) Acenaphthene	14.41	153	9395	0.39	ng/ul	100
9) Fluorene	15.40	166	11524	0.40	ng/ul	100
11) Pentachlorophenol	16.76	266	487	0.16	ng/ul	100
12) Phenanthrene	17.13	178	18853	0.43	ng/ul	100
13) Anthracene	17.23	178	16400	0.41	ng/ul	100
15) Fluoranthene	19.15	202	21336	0.43	ng/ul	100
17) Pyrene	19.53	202	21941	0.43	ng/ul	100
18) Benzo(a)anthracene	21.27	228	16217	0.39	ng/ul	100
19) Chrysene	21.32	228	19240	0.42	ng/ul	100
21) Benzo(b)fluoranthene	22.83	252	16184	0.39	ng/ul	100
22) Benzo(k)fluoranthene	22.89	252	17848	0.45	ng/ul	100
23) Benzo(a)pyrene	23.41	252	15375	0.41	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.75	276	16211	0.39	ng/ul	100
25) Dibenzo(a,h)anthracene	25.76	278	12472	0.38	ng/ul	100
26) Benzo(g,h,i)perylene	26.44	276	14400	0.39	ng/ul	100

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

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