

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041715\
 Data File : BM001024.D
 Acq On : 16 Apr 2015 19:32
 Operator : TP/IZ
 Sample : SSTD0.429
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.429

Quant Time: Apr 17 06:36:15 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM041715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 17 06:23:20 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1439	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	4895	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	3003	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	4792	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	3922	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	3619	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.97	152	2392	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	4000	0.43	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.42	128	5018	0.44	ng/ul	100
5) 2-Methylnaphthalene	12.04	142	3206	0.44	ng/ul	100
7) Acenaphthylene	13.96	152	4115	0.41	ng/ul	100
8) Acenaphthene	14.30	153	3205	0.40	ng/ul	100
9) Fluorene	15.29	166	3643	0.40	ng/ul	100
11) Pentachlorophenol	16.65	266	123	0.22	ng/ul	100
12) Phenanthrene	17.02	178	5526	0.44	ng/ul	100
13) Anthracene	17.12	178	4923	0.44	ng/ul	100
15) Fluoranthene	19.05	202	5899	0.42	ng/ul	100
17) Pyrene	19.43	202	6140	0.44	ng/ul	100
18) Benzo(a)anthracene	21.18	228	4644	0.43	ng/ul	100
19) Chrysene	21.23	228	5644	0.44	ng/ul	100
21) Benzo(b)fluoranthene	22.71	252	5031	0.42	ng/ul	100
22) Benzo(k)fluoranthene	22.75	252	5603	0.45	ng/ul	100
23) Benzo(a)pyrene	23.26	252	4977	0.45	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.53	276	6037	0.46	ng/ul	100
25) Dibenzo(a,h)anthracene	25.54	278	4977	0.47	ng/ul	100
26) Benzo(g,h,i)perylene	26.19	276	5366	0.46	ng/ul	100

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

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