

Data Path : Z:\HPCHEM1\BNA_M\Data\BM111516\
 Data File : BM007910.D
 Acq On : 16 Nov 2016 01:27
 Operator : UM/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.463

Quant Time: Nov 16 07:46:19 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM102516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 16 07:43:19 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	713	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	2338	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	1220	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	2493	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	2439	0.40	ng/ul	0.00
20) Perylene-d12	23.57	264	2319	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.13	152	1177	0.43	ng/ul	0.01
14) Fluoranthene-d10	19.16	212	2553	0.45	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.57	128	2513	0.42	ng/ul	97
5) 2-Methylnaphthalene	12.21	142	1593	0.42	ng/ul	98
7) Acenaphthylene	14.09	152	2299	0.40	ng/ul	98
8) Acenaphthene	14.43	153	1669	0.38	ng/ul	99
9) Fluorene	15.44	166	1867	0.38	ng/ul	100
11) Pentachlorophenol	16.82	266	282	0.48	ng/ul	97
12) Phenanthrene	17.16	178	2835	0.41	ng/ul	98
13) Anthracene	17.26	178	2704	0.43	ng/ul	100
15) Fluoranthene	19.19	202	3577	0.43	ng/ul	97
17) Pyrene	19.56	202	3604	0.43	ng/ul	98
18) Benzo(a)anthracene	21.30	228	3019	0.43	ng/ul	97
19) Chrysene	21.35	228	3378	0.43	ng/ul	93
21) Benzo(b)fluoranthene	22.89	252	3376	0.42	ng/ul	98
22) Benzo(k)fluoranthene	22.93	252	3507	0.42	ng/ul	99
23) Benzo(a)pyrene	23.47	252	3323	0.45	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.84	276	3994	0.45	ng/ul	97
25) Dibenzo(a,h)anthracene	25.85	278	3158	0.44	ng/ul	97
26) Benzo(g,h,i)perylene	26.54	276	3550	0.44	ng/ul	96

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

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