

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122216\
 Data File : BM008582.D
 Acq On : 23 Dec 2016 07:05
 Operator : UM/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.460

Manual Integrations
 APPROVED

umangi
 12/23/2016 6:12:19 PM

Quant Time: Dec 23 07:41:45 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 23 05:50:04 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	309	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	1164	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	660	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	1266	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	1306	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	1292	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	665	0.37	ng/ul	0.01
14) Fluoranthene-d10	19.11	212	1416	0.41	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.54	128	1272	0.39	ng/ul#	77
5) 2-Methylnaphthalene	12.17	142	816	0.37	ng/ul	95
7) Acenaphthylene	14.04	152	1200	0.40	ng/ul#	88
8) Acenaphthene	14.38	153	907	0.40	ng/ul	96
9) Fluorene	15.39	166	998	0.39	ng/ul#	87
11) Pentachlorophenol	16.78	266	160	0.41	ng/ul	93
12) Phenanthrene	17.12	178	1535	0.40	ng/ul#	92
13) Anthracene	17.23	178	1482	0.39	ng/ul#	91
15) Fluoranthene	19.14	202	2011m	0.42	ng/ul	
17) Pyrene	19.50	202	2088	0.41	ng/ul	96
18) Benzo(a)anthracene	21.26	228	1666	0.39	ng/ul	94
19) Chrysene	21.31	228	1932	0.41	ng/ul	94
21) Benzo(b)fluoranthene	22.83	252	1869	0.37	ng/ul	94
22) Benzo(k)fluoranthene	22.87	252	2010	0.40	ng/ul	93
23) Benzo(a)pyrene	23.40	252	1903	0.40	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.75	276	2260	0.40	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.75	278	1794	0.40	ng/ul#	90
26) Benzo(g,h,i)perylene	26.44	276	1943	0.40	ng/ul	94

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

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