

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101215\
 Data File : BE091001.D
 Acq On : 12 Oct 2015 19:17
 Operator : UM/IZ
 Sample : SSTD0.268
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.268

Manual Integrations
 APPROVED

MMdadoda
 10/13/2015 1:26:13 PM

Quant Time: Oct 13 07:18:07 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM02.2-EPA-SIM-BE101215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 13 07:04:09 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	4512	0.40	ng/ul	0.00
2) Naphthalene-d8	9.63	136	20592	0.40	ng/ul	0.00
6) Acenaphthene-d10	13.46	164	11405	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.18	188	29618	0.40	ng/ul	0.01
16) Chrysene-d12	20.33	240	36842	0.40	ng/ul	0.00
20) Perylene-d12	22.24	264	37215	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	13.20	152	41993	0.62	ng/ul	0.00
14) Fluoranthene-d10	18.17	212	15548	0.21	ng/ul	0.00
Target Compounds						Ovalue
3) Naphthalene	9.68	128	10156	0.19	ng/ul	98
5) 2-Methylnaphthalene	11.20	142	6548	0.18	ng/ul	96
7) Acenaphthylene	13.20	152	41993	0.53	ng/ul	99
8) Acenaphthene	13.51	153	30442	0.46	ng/ul	99
9) Fluorene	14.50	166	9400	0.19	ng/ul	98
11) Pentachlorophenol	15.93	266	834	0.38	ng/ul	96
12) Phenanthrene	16.21	178	16593	0.18	ng/ul	100
13) Anthracene	16.29	178	15552	0.18	ng/ul	98
15) Fluoranthene	18.20	202	19406	0.19	ng/ul	99
17) Pyrene	18.59	202	19848	0.14	ng/ul	98
18) Benzo(a)anthracene	20.30	228	19784	0.17	ng/ul	99
19) Chrysene	20.36	228	20917	0.18	ng/ul	99
21) Benzo(b)fluoranthene	21.67	252	22853	0.17	ng/ul	92
22) Benzo(k)fluoranthene	21.70	252	21731	0.17	ng/ul	93
23) Benzo(a)pyrene	22.15	252	21320	0.18	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	23.91	276	22465m	0.18	ng/ul	
25) Dibenzo(a,h)anthracene	23.87	278	22069	0.21	ng/ul	98
26) Benzo(g,h,i)perylene	24.48	276	23604	0.21	ng/ul	96

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

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