

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101215\
 Data File : BE091003.D
 Acq On : 12 Oct 2015 20:29
 Operator : UM/IZ
 Sample : SST0.870
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.870

Manual Integrations
 APPROVED

MMdadoda
 10/13/2015 1:26:14 PM

Quant Time: Oct 13 07:13:43 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	6.96	152	4499	0.40	ng/ul	0.00
2) Naphthalene-d8	9.63	136	20220	0.40	ng/ul	0.00
6) Acenaphthene-d10	13.46	164	10887	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.18	188	26479	0.40	ng/ul	0.01
16) Chrysene-d12	20.33	240	33264	0.40	ng/ul	0.00
20) Perylene-d12	22.24	264	33809	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	13.20	152	171217	2.57	ng/ul	0.00
14) Fluoranthene-d10	18.17	212	61957	0.93	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	9.68	128	42017	0.79	ng/ul	99
5) 2-Methylnaphthalene	11.20	142	27367m	0.76	ng/ul	
7) Acenaphthylene	13.20	152	171217	2.28	ng/ul	100
8) Acenaphthene	13.51	153	123237	1.96	ng/ul	98
9) Fluorene	14.50	166	37720	0.81	ng/ul	99
11) Pentachlorophenol	15.93	266	2605	1.32	ng/ul	94
12) Phenanthrene	16.21	178	63086	0.79	ng/ul	100
13) Anthracene	16.29	178	61097	0.79	ng/ul	100
15) Fluoranthene	18.19	202	76048	0.83	ng/ul	94
17) Pyrene	18.58	202	76562	0.62	ng/ul	99
18) Benzo(a)anthracene	20.30	228	75415	0.70	ng/ul	99
19) Chrysene	20.36	228	80388	0.75	ng/ul	99
21) Benzo(b)fluoranthene	21.67	252	89436	0.74	ng/ul	97
22) Benzo(k)fluoranthene	21.70	252	82782	0.70	ng/ul	98
23) Benzo(a)pyrene	22.14	252	81747	0.74	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	23.92	276	87328	0.77	ng/ul	99
25) Dibenzo(a,h)anthracene	23.87	278	87517	0.93	ng/ul	99
26) Benzo(a,h,i)perylene	24.49	276	92898	0.91	ng/ul	97

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

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