

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080118\
 Data File : BE097129.D
 Acq On : 1 Aug 2018 14:18
 Operator : SJ/JU
 Sample : SSTD0.876
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.876

Quant Time: Aug 01 14:57:19 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 01 14:22:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	1419	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6870	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3738	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10049	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	12672	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	12290	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	8696	0.77	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	26662	1.07	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.18	128	12895	0.776	ng/ul#	91
5) 2-Methylnaphthalene	11.81	142	8701	0.769	ng/ul	98
7) Acenaphthylene	13.73	152	12846	0.838	ng/ul	96
8) Acenaphthene	14.09	153	9280	0.734	ng/ul	95
9) Fluorene	15.08	166	10589	0.811	ng/ul#	92
11) Pentachlorophenol	16.44	266	1088	1.654	ng/ul	98
12) Phenanthrene	16.81	178	19934	0.769	ng/ul#	89
13) Anthracene	16.90	178	18922	0.860	ng/ul#	91
15) Fluoranthene	18.84	202	26249	0.951	ng/ul	84
17) Pyrene	19.20	202	26410	0.635	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	28373	0.949	ng/ul#	85
19) Chrysene	21.01	228	29566	0.764	ng/ul	97
21) Benzo(b)fluoranthene	22.53	252	26932	0.803	ng/ul	82
22) Benzo(k)fluoranthene	22.58	252	28535	0.779	ng/ul#	85
23) Benzo(a)pyrene	23.11	252	26402	0.931	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.50	276	34097	1.158	ng/ul#	76
25) Dibenzo(a,h)anthracene	25.52	278	27915	1.193	ng/ul#	87
26) Benzo(g,h,i)perylene	26.20	276	29676	1.038	ng/ul#	92

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

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