

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080218\
 Data File : BE097144.D
 Acq On : 2 Aug 2018 10:33
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.482

Quant Time: Aug 02 14:47:43 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 : Thu Aug 02 14:47:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1354	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6417	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3551	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	9493	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	11481	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	11425	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	4139	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	12342	0.38	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.18	128	6203	0.399	ng/ul#	93
5) 2-Methylnaphthalene	11.80	142	4099	0.399	ng/ul	97
7) Acenaphthylene	13.73	152	6087	0.396	ng/ul	98
8) Acenaphthene	14.08	153	4434	0.399	ng/ul#	92
9) Fluorene	15.08	166	4973	0.390	ng/ul#	92
11) Pentachlorophenol	16.44	266	456	0.335	ng/ul	98
12) Phenanthrene	16.81	178	9555	0.395	ng/ul#	88
13) Anthracene	16.90	178	8679	0.384	ng/ul#	91
15) Fluoranthene	18.84	202	12389	0.387	ng/ul	84
17) Pyrene	19.20	202	12636	0.411	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	12568	0.383	ng/ul#	85
19) Chrysene	21.01	228	13490	0.399	ng/ul	97
21) Benzo(b)fluoranthene	22.54	252	12188	0.391	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	13048	0.389	ng/ul#	88
23) Benzo(a)pyrene	23.11	252	11876	0.381	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.50	276	14972	0.384	ng/ul#	78
25) Dibenzo(a,h)anthracene	25.52	278	12500	0.392	ng/ul#	89
26) Benzo(g,h,i)perylene	26.20	276	12904	0.383	ng/ul#	92

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

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