

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080118\
 Data File : BE097142.D
 Acq On : 1 Aug 2018 23:08
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.481

Quant Time: Aug 02 01:04:35 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 16:16:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1384	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6627	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3784	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10225	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	12389	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	12217	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	4357	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	13623	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	6439	0.401	ng/ul#	93
5) 2-Methylnaphthalene	11.81	142	4291	0.405	ng/ul	96
7) Acenaphthylene	13.73	152	6484	0.396	ng/ul	97
8) Acenaphthene	14.08	153	4685	0.396	ng/ul#	92
9) Fluorene	15.08	166	5350	0.393	ng/ul	90
11) Pentachlorophenol	16.44	266	570	0.389	ng/ul	96
12) Phenanthrene	16.81	178	10405	0.399	ng/ul#	88
13) Anthracene	16.90	178	9741	0.401	ng/ul#	91
15) Fluoranthene	18.84	202	13364	0.388	ng/ul	83
17) Pyrene	19.20	202	13869	0.418	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	13825	0.391	ng/ul#	85
19) Chrysene	21.01	228	14526	0.398	ng/ul	97
21) Benzo(b)fluoranthene	22.54	252	13137	0.395	ng/ul	84
22) Benzo(k)fluoranthene	22.58	252	14054	0.391	ng/ul#	88
23) Benzo(a)pyrene	23.11	252	12749	0.383	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.50	276	15674	0.376	ng/ul#	77
25) Dibenzo(a,h)anthracene	25.52	278	13244	0.388	ng/ul#	87
26) Benzo(g,h,i)perylene	26.20	276	13493	0.374	ng/ul#	90

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

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