

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102618\
 Data File : BE098081.D
 Acq On : 26 Oct 2018 9:59
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.471

Quant Time: Oct 26 23:14:53 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 26 07:43:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	1014	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	4873	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3420	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8451	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	11491	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	9697	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	3413	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	10165	0.37	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	4504	0.386	ng/ul	98
5) 2-Methylnaphthalene	12.35	142	3415	0.387	ng/ul	99
7) Acenaphthylene	14.25	152	5181	0.384	ng/ul	99
8) Acenaphthene	14.60	153	3799	0.377	ng/ul	100
9) Fluorene	15.59	166	4586	0.378	ng/ul	100
11) Pentachlorophenol	16.95	266	129	0.292	ng/ul	90
12) Phenanthrene	17.33	178	7339	0.376	ng/ul	99
13) Anthracene	17.42	178	7246	0.376	ng/ul	99
15) Fluoranthene	19.35	202	9194	0.370	ng/ul	99
17) Pyrene	19.71	202	10717	0.399	ng/ul	99
18) Benzo(a)anthracene	21.45	228	12354	0.366	ng/ul	100
19) Chrysene	21.51	228	9949	0.381	ng/ul	100
21) Benzo(b)fluoranthene	23.24	252	10046	0.374	ng/ul	96
22) Benzo(k)fluoranthene	23.29	252	9906	0.377	ng/ul	99
23) Benzo(a)pyrene	23.91	252	9753	0.377	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.71	276	11255	0.380	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.74	278	9383	0.380	ng/ul	99
26) Benzo(g,h,i)perylene	27.55	276	9390	0.377	ng/ul	99

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

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