

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101018\
 Data File : BE097820.D
 Acq On : 10 Oct 2018 9:45
 Operator : SJ/JU
 Sample : SSTD0.236
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.236

Quant Time: Oct 10 11:13:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE101018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 10 11:08:35 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2903	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	12482	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	7179	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	20262	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	32140	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	28417	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	3691	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	12411	0.19	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.74	128	5911	0.194	ng/ul	97
5) 2-Methylnaphthalene	12.36	142	3698	0.192	ng/ul	99
7) Acenaphthylene	14.26	152	4885	0.187	ng/ul#	94
8) Acenaphthene	14.60	153	4110	0.183	ng/ul	99
9) Fluorene	15.60	166	4798	0.186	ng/ul	98
11) Pentachlorophenol	16.94	266	350	0.175	ng/ul	96
12) Phenanthrene	17.34	178	9250	0.177	ng/ul	96
13) Anthracene	17.43	178	7489	0.171	ng/ul	97
15) Fluoranthene	19.36	202	11709	0.175	ng/ul	98
17) Pyrene	19.72	202	14836	0.172	ng/ul	98
18) Benzo(a)anthracene	21.46	228	15271	0.185	ng/ul	98
19) Chrysene	21.52	228	16080	0.169	ng/ul	97
21) Benzo(b)fluoranthene	23.25	252	14012	0.184	ng/ul	97
22) Benzo(k)fluoranthene	23.29	252	14966	0.176	ng/ul	95
23) Benzo(a)pyrene	23.91	252	12341	0.175	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.71	276	13832	0.152	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.74	278	11348	0.147	ng/ul	96
26) Benzo(g,h,i)perylene	27.53	276	13329	0.171	ng/ul	99

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

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