

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE103018\
 Data File : BE098108.D
 Acq On : 30 Oct 2018 10:57
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.474

Quant Time: Oct 30 12:54:28 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 : Tue Oct 30 01:14:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	1072	0.40	ng/ul	0.00
2) Naphthalene-d8	10.67	136	5221	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3631	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8818	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	13240	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	11005	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	3503	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	10557	0.37	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.73	128	4735	0.379	ng/ul	97
5) 2-Methylnaphthalene	12.34	142	3557	0.376	ng/ul	100
7) Acenaphthylene	14.25	152	5398	0.377	ng/ul	98
8) Acenaphthene	14.59	153	3956	0.370	ng/ul	97
9) Fluorene	15.59	166	4739	0.368	ng/ul	98
11) Pentachlorophenol	16.95	266	223	0.483	ng/ul	98
12) Phenanthrene	17.33	178	7628	0.375	ng/ul	99
13) Anthracene	17.42	178	7490	0.373	ng/ul	99
15) Fluoranthene	19.35	202	9641	0.372	ng/ul	99
17) Pyrene	19.71	202	9793	0.317	ng/ul	98
18) Benzo(a)anthracene	21.45	228	12702	0.326	ng/ul	100
19) Chrysene	21.51	228	10476	0.348	ng/ul	98
21) Benzo(b)fluoranthene	23.24	252	10375	0.340	ng/ul	97
22) Benzo(k)fluoranthene	23.29	252	10598	0.355	ng/ul	98
23) Benzo(a)pyrene	23.91	252	10198	0.348	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.71	276	11606	0.345	ng/ul#	96
25) Dibenzo(a,h)anthracene	26.74	278	9557	0.341	ng/ul	99
26) Benzo(g,h,i)perylene	27.54	276	9748	0.345	ng/ul	97

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

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