

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102518\
 Data File : BE098052.D
 Acq On : 25 Oct 2018 12:23
 Operator : SJ/JU
 Sample : SSTD0.463
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.463

Quant Time: Oct 25 13:28:13 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 : Thu Oct 25 13:24:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	933	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	4711	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3386	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8316	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	11722	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	9807	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	3326	0.45	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	10784	0.40	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	4536	0.400	ng/ul	100
5) 2-Methylnaphthalene	12.34	142	3408	0.452	ng/ul	100
7) Acenaphthylene	14.25	152	5315	0.446	ng/ul	100
8) Acenaphthene	14.60	153	3953	0.398	ng/ul	100
9) Fluorene	15.59	166	4738	0.395	ng/ul	100
11) Pentachlorophenol	16.95	266	170	0.224	ng/ul	100
12) Phenanthrene	17.33	178	7662	0.386	ng/ul	100
13) Anthracene	17.42	178	7530	0.446	ng/ul	100
15) Fluoranthene	19.35	202	9750	0.372	ng/ul	100
17) Pyrene	19.71	202	11268	0.391	ng/ul	100
18) Benzo(a)anthracene	21.45	228	14287	0.439	ng/ul	100
19) Chrysene	21.51	228	10705	0.341	ng/ul	100
21) Benzo(b)fluoranthene	23.24	252	10850	0.428	ng/ul	100
22) Benzo(k)fluoranthene	23.29	252	10440	0.363	ng/ul	100
23) Benzo(a)pyrene	23.91	252	10259	0.442	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.71	276	11667	0.443	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.74	278	9859	0.448	ng/ul	100
26) Benzo(g,h,i)perylene	27.55	276	10038	0.404	ng/ul	100

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

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