

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE103118\
 Data File : BE098119.D
 Acq On : 31 Oct 2018 16:59
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.477

Quant Time: Nov 01 01:45:31 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 Nov 01 01:42:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	1078	0.40	ng/ul	0.00
2) Naphthalene-d8	10.67	136	5240	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3466	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.28	188	8433	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	11769	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	9687	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	3646	0.39	ng/ul	-0.01
14) Fluoranthene-d10	19.32	212	10761	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.72	128	4986	0.398	ng/ul	97
5) 2-Methylnaphthalene	12.34	142	3696	0.390	ng/ul	100
7) Acenaphthylene	14.24	152	5506	0.403	ng/ul	98
8) Acenaphthene	14.59	153	4055	0.397	ng/ul	98
9) Fluorene	15.58	166	4818	0.392	ng/ul	98
11) Pentachlorophenol	16.95	266	235	0.533	ng/ul	93
12) Phenanthrene	17.33	178	7747	0.398	ng/ul	98
13) Anthracene	17.42	178	7603	0.396	ng/ul	99
15) Fluoranthene	19.35	202	9864	0.398	ng/ul	97
17) Pyrene	19.71	202	11752	0.428	ng/ul	97
18) Benzo(a)anthracene	21.44	228	12192	0.352	ng/ul	99
19) Chrysene	21.50	228	10488	0.392	ng/ul	100
21) Benzo(b)fluoranthene	23.23	252	10756	0.401	ng/ul	99
22) Benzo(k)fluoranthene	23.28	252	10349	0.394	ng/ul	99
23) Benzo(a)pyrene	23.90	252	10235	0.397	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.71	276	11725	0.396	ng/ul#	95
25) Dibenzo(a,h)anthracene	26.73	278	9799	0.397	ng/ul	97
26) Benzo(g,h,i)perylene	27.53	276	9772	0.392	ng/ul	96

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

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