

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120718\
 Data File : BN003773.D
 Acq On : 07 Dec 2018 19:04
 Operator : JU/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.493

Quant Time: Dec 08 00:37:50 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 08 00:32:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	5758	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	23384	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	12063	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	26003	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	22174	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	18518	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	14388	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	29018	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	27268	0.415	ng/ul	99
5) 2-Methylnaphthalene	12.29	142	18937	0.413	ng/ul	99
7) Acenaphthylene	14.19	152	23444	0.432	ng/ul	100
8) Acenaphthene	14.53	153	19818	0.416	ng/ul	99
9) Fluorene	15.52	166	21991	0.396	ng/ul	99
11) Pentachlorophenol	16.86	266	2152	0.459	ng/ul	95
12) Phenanthrene	17.26	178	36001	0.422	ng/ul	100
13) Anthracene	17.34	178	30477	0.435	ng/ul	100
15) Fluoranthene	19.27	202	39927	0.396	ng/ul	97
17) Pyrene	19.63	202	38921	0.489	ng/ul	99
18) Benzo(a)anthracene	21.38	228	31862	0.455	ng/ul	99
19) Chrysene	21.43	228	32989	0.412	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	31782	0.415	ng/ul	99
22) Benzo(k)fluoranthene	23.05	252	30972	0.415	ng/ul	99
23) Benzo(a)pyrene	23.61	252	28160	0.425	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.05	276	30958	0.407	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.07	278	24665	0.403	ng/ul	99
26) Benzo(g,h,i)perylene	26.78	276	25935	0.392	ng/ul	99

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

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