

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120718\
 Data File : BN003758.D
 Acq On : 07 Dec 2018 10:10
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.492

Quant Time: Dec 07 10:57:38 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 : Fri Dec 07 04:08:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	5935	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	24254	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	12472	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	27321	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	24966	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	21566	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	14862	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	33803	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	28221	0.414	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	19629	0.413	ng/ul	100
7) Acenaphthylene	14.19	152	24338	0.433	ng/ul	100
8) Acenaphthene	14.53	153	20694	0.421	ng/ul	99
9) Fluorene	15.52	166	22971	0.400	ng/ul	99
11) Pentachlorophenol	16.86	266	2231	0.453	ng/ul	100
12) Phenanthrene	17.26	178	37607	0.419	ng/ul	100
13) Anthracene	17.34	178	32093	0.435	ng/ul	100
15) Fluoranthene	19.27	202	45386	0.429	ng/ul	98
17) Pyrene	19.63	202	43393	0.485	ng/ul	97
18) Benzo(a)anthracene	21.38	228	37262	0.473	ng/ul	99
19) Chrysene	21.43	228	37013	0.410	ng/ul	100
21) Benzo(b)fluoranthene	23.00	252	35749	0.401	ng/ul	99
22) Benzo(k)fluoranthene	23.05	252	36056	0.415	ng/ul	99
23) Benzo(a)pyrene	23.60	252	32285	0.418	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.05	276	34712	0.392	ng/ul	98
25) Dibenzo(a,h)anthracene	26.06	278	27612	0.387	ng/ul	99
26) Benzo(g,h,i)perylene	26.77	276	29052	0.377	ng/ul	98

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

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