

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101119\
 Data File : BN008109.D
 Acq On : 12 Oct 2019 12:21
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.412

Quant Time: Oct 12 13:30:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 11 19:05:54 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2279	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	11542	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	7566	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	18873	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	19793	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	18574	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	7471	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	24745	0.38	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.45	128	11593	0.353	ng/ul	99
5) 2-Methylnaphthalene	12.07	142	8423	0.374	ng/ul	100
7) Acenaphthylene	13.98	152	14406	0.400	ng/ul	100
8) Acenaphthene	14.32	153	11403	0.407	ng/ul	98
9) Fluorene	15.32	166	11750	0.369	ng/ul	97
11) Pentachlorophenol	16.68	266	1691	0.370	ng/ul	98
12) Phenanthrene	17.05	178	19825	0.359	ng/ul	100
13) Anthracene	17.14	178	18875	0.388	ng/ul	100
15) Fluoranthene	19.07	202	26557	0.364	ng/ul	98
17) Pyrene	19.44	202	30828	0.391	ng/ul	99
18) Benzo(a)anthracene	21.20	228	26812	0.373	ng/ul	97
19) Chrysene	21.25	228	29724	0.337	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	25426	0.421	ng/ul	90
22) Benzo(k)fluoranthene	22.79	252	26814	0.391	ng/ul#	92
23) Benzo(a)pyrene	23.31	252	24588	0.386	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.59	276	24497	0.327	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.60	278	19906	0.336	ng/ul#	88
26) Benzo(g,h,i)perylene	26.25	276	20717	0.301	ng/ul#	92

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

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