

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101619\
 Data File : BN008202.D
 Acq On : 16 Oct 2019 17:17
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.421

Quant Time: Oct 16 18:30:03 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 : Wed Oct 16 18:17:24 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	2387	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	11702	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	7328	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	16470	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	15009	0.40	ng/ul	0.00
20) Perylene-d12	23.38	264	12407	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	7528	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	20572	0.36	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.42	128	12592	0.378	ng/ul	99
5) 2-Methylnaphthalene	12.05	142	8781	0.384	ng/ul	99
7) Acenaphthylene	13.96	152	13001	0.372	ng/ul	100
8) Acenaphthene	14.31	153	11029	0.406	ng/ul	99
9) Fluorene	15.30	166	11703	0.380	ng/ul	98
11) Pentachlorophenol	16.66	266	1446	0.363	ng/ul	99
12) Phenanthrene	17.03	178	18949	0.393	ng/ul	99
13) Anthracene	17.13	178	16912	0.399	ng/ul	100
15) Fluoranthene	19.06	202	23279	0.366	ng/ul	97
17) Pyrene	19.42	202	25828	0.432	ng/ul	97
18) Benzo(a)anthracene	21.18	228	21377	0.392	ng/ul	99
19) Chrysene	21.23	228	23898	0.357	ng/ul	100
21) Benzo(b)fluoranthene	22.73	252	19001	0.470	ng/ul	94
22) Benzo(k)fluoranthene	22.77	252	19941	0.436	ng/ul	98
23) Benzo(a)pyrene	23.28	252	17544	0.413	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.54	276	17073	0.341	ng/ul	97
25) Dibenzo(a,h)anthracene	25.56	278	13701	0.346	ng/ul	99
26) Benzo(g,h,i)perylene	26.20	276	15058	0.328	ng/ul	98

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

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