

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101119\
 Data File : BN008090.D
 Acq On : 11 Oct 2019 22:40
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.410

Quant Time: Oct 11 23:16:57 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	2062	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	9887	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	6289	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	14958	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	15356	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	17070	0.40	ng/ul	-0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	5960	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	18304	0.35	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.45	128	10220	0.363	ng/ul	99
5) 2-Methylnaphthalene	12.07	142	6829	0.354	ng/ul	100
7) Acenaphthylene	13.98	152	10607	0.354	ng/ul	99
8) Acenaphthene	14.32	153	9364	0.402	ng/ul	99
9) Fluorene	15.32	166	9578	0.362	ng/ul	98
11) Pentachlorophenol	16.67	266	1259	0.348	ng/ul	99
12) Phenanthrene	17.05	178	15339	0.350	ng/ul	99
13) Anthracene	17.14	178	13184	0.342	ng/ul	100
15) Fluoranthene	19.07	202	20029	0.346	ng/ul	100
17) Pyrene	19.44	202	23509	0.384	ng/ul	99
18) Benzo(a)anthracene	21.19	228	18966	0.340	ng/ul	99
19) Chrysene	21.24	228	25220	0.368	ng/ul	99
21) Benzo(b)fluoranthene	22.74	252	21942	0.395	ng/ul	99
22) Benzo(k)fluoranthene	22.78	252	24927	0.396	ng/ul	99
23) Benzo(a)pyrene	23.30	252	23039	0.394	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.57	276	27602	0.401	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.58	278	21697	0.399	ng/ul	99
26) Benzo(g,h,i)perylene	26.23	276	25340	0.401	ng/ul	99

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

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