

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101019\
 Data File : BN008062.D
 Acq On : 10 Oct 2019 14:40
 Operator : JU
 Sample : SSTD1.605
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.605

Quant Time: Oct 10 16:14:38 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 : Thu Oct 10 16:00:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2550	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	12404	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	7626	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	18582	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	18982	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	21917	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	33987	1.76	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	100802	1.67	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.45	128	54560	1.558	ng/ul	98
5) 2-Methylnaphthalene	12.07	142	39331	1.672	ng/ul	99
7) Acenaphthylene	13.99	152	60213	1.419	ng/ul	98
8) Acenaphthene	14.33	153	45297	1.486	ng/ul	99
9) Fluorene	15.32	166	53422	1.619	ng/ul	98
11) Pentachlorophenol	16.68	266	7598	1.815	ng/ul	99
12) Phenanthrene	17.06	178	87136	1.503	ng/ul	99
13) Anthracene	17.15	178	81978	1.500	ng/ul	98
15) Fluoranthene	19.08	202	113062	1.457	ng/ul	99
17) Pyrene	19.44	202	116304	1.252	ng/ul	98
18) Benzo(a)anthracene	21.20	228	114622	1.351	ng/ul	99
19) Chrysene	21.25	228	125856	1.381	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	121273	1.267	ng/ul	90
22) Benzo(k)fluoranthene	22.80	252	127101	1.270	ng/ul#	90
23) Benzo(a)pyrene	23.31	252	120889	1.267	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.59	276	144635	1.354	ng/ul	99
25) Dibenzo(a,h)anthracene	25.60	278	116577	1.379	ng/ul#	92
26) Benzo(g,h,i)perylene	26.25	276	128742	1.412	ng/ul	96

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

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