

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080719\
 Data File : BN006984.D
 Acq On : 07 Aug 2019 13:06
 Operator : HP/JU
 Sample : SSTD1.623
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.624

Quant Time: Aug 07 17:26:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN080719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 17:13:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	4909	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	19020	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	11972	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	30241	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	28111	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	30118	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	54055	1.60	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	157381	1.61	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.28	128	87589	1.600	ng/ul	97
5) 2-Methylnaphthalene	11.91	142	65584	1.596	ng/ul	100
7) Acenaphthylene	13.83	152	111267	1.604	ng/ul	99
8) Acenaphthene	14.18	153	75806	1.597	ng/ul	99
9) Fluorene	15.17	166	91445	1.589	ng/ul	100
11) Pentachlorophenol	16.52	266	13210	1.252	ng/ul#	80
12) Phenanthrene	16.90	178	153769	1.596	ng/ul	99
13) Anthracene	16.99	178	157078	1.593	ng/ul	99
15) Fluoranthene	18.93	202	191745	1.601	ng/ul	99
17) Pyrene	19.30	202	195620	1.559	ng/ul	98
18) Benzo(a)anthracene	21.07	228	198571	1.608	ng/ul	99
19) Chrysene	21.12	228	184055	1.592	ng/ul	99
21) Benzo(b)fluoranthene	22.58	252	197268	1.616	ng/ul	98
22) Benzo(k)fluoranthene	22.63	252	187002	1.580	ng/ul	97
23) Benzo(a)pyrene	23.12	252	185459	1.613	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.31	276	205227	1.656	ng/ul	99
25) Dibenzo(a,h)anthracene	25.33	278	163801	1.655	ng/ul	97
26) Benzo(g,h,i)perylene	25.94	276	171527	1.650	ng/ul	97

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

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