

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN020421\
 Data File : BN013568.D
 Acq On : 04 Feb 2021 13:07
 Operator : CG/JU
 Sample : SSTD0.805
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.805

Quant Time: Feb 04 13:36:31 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN020421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 04 13:01:49 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	4536	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	16414	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	8694	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	18235	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	17080	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	13047	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	18363	0.75	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	35459	0.73	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	33269	0.701	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	22254	0.738	ng/ul	100
7) Acenaphthylene	13.91	152	25290	0.706	ng/ul	99
8) Acenaphthene	14.25	153	22470	0.705	ng/ul	99
9) Fluorene	15.25	166	25394	0.720	ng/ul	99
11) Pentachlorophenol	16.59	266	1613	0.600	ng/ul	100
12) Phenanthrene	16.98	178	41831	0.701	ng/ul	99
13) Anthracene	17.07	178	35176	0.718	ng/ul	100
15) Fluoranthene	19.00	202	45353	0.695	ng/ul	98
17) Pyrene	19.37	202	44989	0.644	ng/ul	100
18) Benzo(a)anthracene	21.12	228	42568	0.755	ng/ul	99
19) Chrysene	21.18	228	46214	0.657	ng/ul	100
21) Benzo(b)fluoranthene	22.66	252	40801	0.745	ng/ul	99
22) Benzo(k)fluoranthene	22.71	252	43899	0.753	ng/ul	97
23) Benzo(a)pyrene	23.21	252	33307	0.717	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.46	276	44952	0.728	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.47	278	37142	0.745	ng/ul	98
26) Benzo(g,h,i)perylene	26.11	276	39133	0.716	ng/ul	100

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

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