

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121619\
 Data File : BN008995.D
 Acq On : 16 Dec 2019 12:47
 Operator : JU
 Sample : SSTD0.404
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.404

Quant Time: Dec 16 13:37:04 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 16 13:30:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	2710	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	11327	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	7716	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.92	188	17333	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	12873	0.40	ng/ul	0.00
20) Perylene-d12	23.27	264	10913	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	8313	0.45	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	20642	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	14050	0.411	ng/ul	100
5) 2-Methylnaphthalene	11.97	142	10672	0.448	ng/ul	100
7) Acenaphthylene	13.89	152	15768	0.370	ng/ul	100
8) Acenaphthene	14.24	153	12347	0.390	ng/ul	100
9) Fluorene	15.23	166	14562	0.392	ng/ul	100
11) Pentachlorophenol	16.59	266	1627	0.248	ng/ul	100
12) Phenanthrene	16.96	178	24600	0.421	ng/ul	100
13) Anthracene	17.05	178	21900	0.396	ng/ul	100
15) Fluoranthene	18.99	202	27748	0.386	ng/ul	100
17) Pyrene	19.35	202	27274	0.528	ng/ul	100
18) Benzo(a)anthracene	21.11	228	23621	0.456	ng/ul	100
19) Chrysene	21.17	228	23943	0.471	ng/ul	100
21) Benzo(b)fluoranthene	22.64	252	20885	0.464	ng/ul	100
22) Benzo(k)fluoranthene	22.68	252	20410	0.457	ng/ul	100
23) Benzo(a)pyrene	23.18	252	18213	0.431	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.38	276	19589	0.397	ng/ul	100
25) Dibenzo(a,h)anthracene	25.39	278	15647	0.392	ng/ul	100
26) Benzo(g,h,i)perylene	26.02	276	17007	0.412	ng/ul	100

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

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