

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN101420\
 Data File : BN012192.D
 Acq On : 14 Oct 2020 13:32
 Operator : CG/JU
 Sample : SST00.406
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.406

Quant Time: Oct 14 14:10:45 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 14 14:09:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1292	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	5219	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3031	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	6079	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	5295	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	5578	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	3094	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	6448	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	5883	0.381	ng/ul	100
5) 2-Methylnaphthalene	12.07	142	3801	0.369	ng/ul	100
7) Acenaphthylene	13.99	152	5028	0.344	ng/ul	100
8) Acenaphthene	14.33	153	4281	0.365	ng/ul	100
9) Fluorene	15.32	166	4743	0.350	ng/ul	100
11) Pentachlorophenol	16.67	266	665	0.362	ng/ul	100
12) Phenanthrene	17.06	178	7523	0.368	ng/ul	100
13) Anthracene	17.15	178	6449	0.354	ng/ul	100
15) Fluoranthene	19.08	202	8729	0.384	ng/ul	100
17) Pyrene	19.45	202	9210	0.392	ng/ul	100
18) Benzo(a)anthracene	21.20	228	7573	0.368	ng/ul	100
19) Chrysene	21.25	228	9306	0.405	ng/ul	100
21) Benzo(b)fluoranthene	22.75	252	8946	0.342	ng/ul	100
22) Benzo(k)fluoranthene	22.80	252	9806	0.366	ng/ul	100
23) Benzo(a)pyrene	23.31	252	8494	0.376	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.59	276	11092	0.374	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.60	278	8880	0.367	ng/ul	100
26) Benzo(g,h,i)perylene	26.26	276	10172	0.392	ng/ul	100

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

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