

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101420\
 Data File : BN012190.D
 Acq On : 14 Oct 2020 12:20
 Operator : CG/JU
 Sample : SSTDO.104
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDO.104

Quant Time: Oct 14 14:13:18 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	1239	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	4840	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	2775	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	5653	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	5025	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	5481	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	658	0.09	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	1538	0.10	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.45	128	1412	0.099	ng/ul#	90
5) 2-Methylnaphthalene	12.08	142	861	0.090	ng/ul	96
7) Acenaphthylene	13.99	152	1160	0.087	ng/ul#	91
8) Acenaphthene	14.33	153	1008	0.094	ng/ul	97
9) Fluorene	15.32	166	1107	0.089	ng/ul#	95
12) Phenanthrene	17.06	178	1792	0.094	ng/ul#	90
13) Anthracene	17.15	178	1412	0.083	ng/ul#	87
15) Fluoranthene	19.08	202	2104	0.100	ng/ul	91
17) Pyrene	19.45	202	2319	0.104	ng/ul#	92
18) Benzo(a)anthracene	21.20	228	1778	0.091	ng/ul#	93
19) Chrysene	21.26	228	2333	0.107	ng/ul	97
21) Benzo(b)fluoranthene	22.76	252	2066	0.080	ng/ul#	48
22) Benzo(k)fluoranthene	22.80	252	2429	0.092	ng/ul#	53
23) Benzo(a)pyrene	23.32	252	2125	0.096	ng/ul#	37
24) Indeno(1,2,3-cd)pyrene	25.60	276	2624	0.090	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.62	278	2090	0.088	ng/ul#	48
26) Benzo(a,h,i)perylene	26.26	276	2506	0.098	ng/ul#	83

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

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