

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN041221\
 Data File : BN014259.D
 Acq On : 12 Apr 2021 13:57
 Operator : CG/JU
 Sample : SSTDO.106
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDO.1006

Quant Time: Apr 12 16:30:34 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN041221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 12 16:13:58 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	6163	0.40	ng/ul	0.00
4) Naphthalene-d8	10.48	136	20716	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.34	164	10716	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.09	188	21450	0.40	ng/ul	0.00
17) Chrysene-d12	21.28	240	18109	0.40	ng/ul	0.00
23) Perylene-d12	23.51	264	15713	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/ul	
6) 2-Methylnaphthalene-d10	12.08	152	3516	0.12	ng/ul	0.00
18) Fluoranthene-d10	19.12	212	5869	0.11	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	10.53	128	6369	0.106	ng/ul	96
7) 2-Methylnaphthalene	12.15	142	4023	0.102	ng/ul	100
8) 1-Methylnaphthalene	12.37	142	3883	0.102	ng/ul	100
10) Acenaphthylene	14.06	152	5146	0.107	ng/ul	97
11) Acenaphthene	14.40	153	4015	0.100	ng/ul	98
12) Fluorene	15.39	166	4492	0.100	ng/ul#	99
15) Phenanthrene	17.13	178	7064	0.098	ng/ul	97
16) Anthracene	17.22	178	6166	0.104	ng/ul	97
19) Fluoranthene	19.15	202	7611	0.105	ng/ul	97
20) Pyrene	19.51	202	7878	0.104	ng/ul	94
21) Benzo(a)anthracene	21.27	228	6928	0.107	ng/ul	98
22) Chrysene	21.32	228	7224	0.101	ng/ul	98
24) Benzo(b)fluoranthene	22.84	252	7212	0.098	ng/ul	82
25) Benzo(k)fluoranthene	22.89	252	7178	0.097	ng/ul#	83
26) Benzo(a)pyrene	23.42	252	5994	0.101	ng/ul#	72
27) Indeno(1,2,3-cd)pyrene	25.76	276	7951	0.100	ng/ul#	48
28) Dibenzo(a,h)anthracene	25.78	278	6349	0.099	ng/ul#	89
29) Benzo(a,h,i)perylene	26.45	276	6866	0.097	ng/ul	91

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

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