

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN041221\
 Data File : BN014264.D
 Acq On : 12 Apr 2021 17:18
 Operator : CG/JU
 Sample : SSTD3.205
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.2005

Quant Time: Apr 12 17:49:10 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	5882	0.40	ng/ul	0.00
4) Naphthalene-d8	10.48	136	20218	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.34	164	10249	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.09	188	20709	0.40	ng/ul	0.00
17) Chrysene-d12	21.28	240	19034	0.40	ng/ul	0.00
23) Perylene-d12	23.51	264	15129	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.23	96	19688	3.12	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.08	152	90854	3.11	ng/ul	0.00
18) Fluoranthene-d10	19.12	212	179107	3.15	ng/ul	0.00
Target Compounds				Ovalue		
2) 1,4-Dioxane	3.27	88	20833	3.159	ng/ul#	73
5) Naphthalene	10.53	128	174635	2.965	ng/ul	99
7) 2-Methylnaphthalene	12.15	142	116435	3.038	ng/ul	100
8) 1-Methylnaphthalene	12.37	142	111176	3.001	ng/ul	100
10) Acenaphthylene	14.06	152	152657	3.304	ng/ul	99
11) Acenaphthene	14.40	153	116846	3.045	ng/ul	98
12) Fluorene	15.39	166	132233	3.086	ng/ul	98
14) Pentachlorophenol	16.73	266	21190	5.984	ng/ul	99
15) Phenanthrene	17.13	178	209192	3.021	ng/ul	99
16) Anthracene	17.22	178	190879	3.343	ng/ul	99
19) Fluoranthene	19.15	202	235302	3.076	ng/ul	98
20) Pyrene	19.51	202	232572	2.933	ng/ul	98
21) Benzo(a)anthracene	21.26	228	219767	3.243	ng/ul	99
22) Chrysene	21.31	228	221136	2.952	ng/ul	99
24) Benzo(b)fluoranthene	22.84	252	217556	3.072	ng/ul	95
25) Benzo(k)fluoranthene	22.89	252	224142	3.132	ng/ul	95
26) Benzo(a)pyrene	23.41	252	184445	3.213	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	25.76	276	253937	3.331	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.78	278	204041	3.320	ng/ul	96
29) Benzo(g,h,i)perylene	26.45	276	212231	3.128	ng/ul	98

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

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