

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN051624\
 Data File : BN031531.D
 Acq On : 15 May 2024 10:40
 Operator : MA/JU
 Sample : SSTD0.488
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4218

Quant Time: May 15 12:16:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN051624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 15 12:15:51 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.700	152	11627	0.400	ng/u1	0.00
4) Naphthalene-d8	10.480	136	36377	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.335	164	19119	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.078	188	42115	0.400	ng/u1	0.00
17) Chrysene-d12	21.266	240	34832	0.400	ng/u1	0.00
23) Perylene-d12	23.502	264	31458	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	5246	0.332	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.074	152	19242	0.299	ng/u1	0.00
18) Fluoranthene-d10	19.109	212	36259	0.288	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.244	88	5257	0.326	ng/u1	95
5) Naphthalene	10.529	128	36528	0.308	ng/u1	99
7) 2-Methylnaphthalene	12.151	142	23035	0.299	ng/u1	100
8) 1-Methylnaphthalene	12.371	142	24285	0.298	ng/u1	100
10) Acenaphthylene	14.053	152	32336	0.309	ng/u1	100
11) Acenaphthene	14.400	153	24574	0.304	ng/u1	99
12) Fluorene	15.385	166	27066	0.289	ng/u1	99
14) Pentachlorophenol	16.723	266	2339	0.263	ng/u1	100
15) Phenanthrene	17.120	178	45415	0.299	ng/u1	100
16) Anthracene	17.213	178	38104	0.295	ng/u1	100
19) Fluoranthene	19.137	202	48695	0.288	ng/u1	100
20) Pyrene	19.499	202	47864	0.288	ng/u1	99
21) Benzo(a)anthracene	21.251	228	42031	0.290	ng/u1	99
22) Chrysene	21.301	228	47537	0.300	ng/u1	99
24) Benzo(b)fluoranthene	22.829	252	41150	0.270	ng/u1	98
25) Benzo(k)fluoranthene	22.873	252	45461	0.281	ng/u1	98
26) Benzo(a)pyrene	23.402	252	35303	0.282	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	25.762	276	44172	0.310	ng/u1	99
28) Dibenzo(a,h)anthracene	25.782	278	34888	0.308	ng/u1	99
29) Benzo(g,h,i)perylene	26.450	276	36140	0.304	ng/u1	100

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

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