

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080322\
 Data File : BN020980.D
 Acq On : 03 Aug 2022 16:21
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV236

Quant Time: Aug 03 16:54:28 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 03 16:31:11 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	3064	0.400	ng/u1	0.00
4) Naphthalene-d8	10.607	136	9617	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.457	164	5375	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.210	188	10518	0.400	ng/u1	0.00
17) Chrysene-d12	21.421	240	8505	0.400	ng/u1	0.00
23) Perylene-d12	23.768	264	6882	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	1444	0.368	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.208	152	5514	0.391	ng/u1	0.00
18) Fluoranthene-d10	19.249	212	10888	0.415	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.245	88	1267	0.314	ng/u1	91
5) Naphthalene	10.651	128	11319	0.383	ng/u1	98
7) 2-Methylnaphthalene	12.279	142	6835	0.381	ng/u1	99
8) 1-Methylnaphthalene	12.499	142	7016	0.391	ng/u1	100
10) Acenaphthylene	14.179	152	9484	0.377	ng/u1	99
11) Acenaphthene	14.521	153	7898	0.380	ng/u1	100
12) Fluorene	15.512	166	8467	0.373	ng/u1	99
14) Pentachlorophenol	16.885	266	498	0.307	ng/u1	98
15) Phenanthrene	17.252	178	13643	0.363	ng/u1	100
16) Anthracene	17.349	178	10393	0.357	ng/u1	100
19) Fluoranthene	19.282	202	14346	0.381	ng/u1	99
20) Pyrene	19.644	202	14889	0.383	ng/u1	95
21) Benzo(a)anthracene	21.410	228	9935	0.358	ng/u1	99
22) Chrysene	21.459	228	13574	0.371	ng/u1	99
24) Benzo(b)fluoranthene	23.055	252	12630	0.396	ng/u1	99
25) Benzo(k)fluoranthene	23.105	252	13330	0.396	ng/u1	98
26) Benzo(a)pyrene	23.663	252	10366	0.357	ng/u1	98
27) Indeno(1,2,3-cd)pyrene	26.189	276	11939	0.344	ng/u1	97
28) Dibenzo(a,h)anthracene	26.209	278	9629	0.351	ng/u1	97
29) Benzo(g,h,i)perylene	26.927	276	11284	0.364	ng/u1	98

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

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