

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031524\
 Data File : BN030409.D
 Acq On : 14 Mar 2024 22:49
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV201

Quant Time: Mar 15 01:05:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN031524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 15 01:04:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.604	152	1867	0.400	ng/ul	0.00
4) Naphthalene-d8	10.377	136	4210	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.254	164	2426	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.013	188	4856	0.400	ng/ul	0.00
17) Chrysene-d12	21.218	240	4417	0.400	ng/ul	0.00
23) Perylene-d12	23.425	264	5159	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.157	96	920	0.404	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.983	152	2327	0.400	ng/ul	0.00
18) Fluoranthene-d10	19.051	212	5344	0.393	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.186	88	916	0.370	ng/ul	92
5) Naphthalene	10.427	128	4687	0.412	ng/ul	99
7) 2-Methylnaphthalene	12.060	142	2764	0.415	ng/ul#	100
8) 1-Methylnaphthalene	12.274	142	2775	0.393	ng/ul#	28
10) Acenaphthylene	13.972	152	4691	0.406	ng/ul	99
11) Acenaphthene	14.319	153	3364	0.397	ng/ul	100
12) Fluorene	15.313	166	3794	0.407	ng/ul	98
14) Pentachlorophenol	16.666	266	512	0.367	ng/ul#	100
15) Phenanthrene	17.055	178	5671	0.396	ng/ul	98
16) Anthracene	17.152	178	5051	0.386	ng/ul	99
19) Fluoranthene	19.079	202	6903	0.395	ng/ul	99
20) Pyrene	19.441	202	7483	0.401	ng/ul	99
21) Benzo(a)anthracene	21.201	228	5802	0.392	ng/ul	99
22) Chrysene	21.253	228	6952	0.408	ng/ul	98
24) Benzo(b)fluoranthene	22.761	252	7045	0.407	ng/ul	96
25) Benzo(k)fluoranthene	22.808	252	8279	0.393	ng/ul	98
26) Benzo(a)pyrene	23.325	252	7169	0.392	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	25.637	276	9615	0.398	ng/ul#	90
28) Dibenzo(a,h)anthracene	25.661	278	7313	0.397	ng/ul	95
29) Benzo(g,h,i)perylene	26.311	276	8277	0.384	ng/ul	98

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

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