

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN012321\
 Data File : BN013470.D
 Acq On : 23 Jan 2021 00:10
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV002

Quant Time: Jan 23 02:39:09 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN012321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 23 02:38:34 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.574	152	3802	0.40	ng/ul	0.00
4) Naphthalene-d8	10.334	136	14041	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.209	164	7245	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.953	188	15322	0.40	ng/ul	0.00
17) Chrysene-d12	21.157	240	15272	0.40	ng/ul	0.00
23) Perylene-d12	23.328	264	13122	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.178	96	1433	0.36	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.929	152	8392	0.42	ng/ul	0.00
18) Fluoranthene-d10	18.990	212	16001	0.40	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.207	88	1557	0.40	ng/ul	98
5) Naphthalene	10.384	128	16606	0.42	ng/ul	100
7) 2-Methylnaphthalene	12.006	142	11044	0.43	ng/ul	99
8) 1-Methylnaphthalene	12.226	142	10345	0.42	ng/ul	100
10) Acenaphthylene	13.922	152	12976	0.43	ng/ul	100
11) Acenaphthene	14.269	153	10919	0.42	ng/ul	100
12) Fluorene	15.263	166	12324	0.42	ng/ul	100
14) Pentachlorophenol	16.611	266	823	0.34	ng/ul	97
15) Phenanthrene	16.996	178	20524	0.42	ng/ul	99
16) Anthracene	17.089	178	16921	0.41	ng/ul	100
19) Fluoranthene	19.023	202	22325	0.40	ng/ul	100
20) Pyrene	19.385	202	22826	0.40	ng/ul	98
21) Benzo(a)anthracene	21.139	228	21870	0.41	ng/ul	100
22) Chrysene	21.192	228	23903	0.43	ng/ul	99
24) Benzo(b)fluoranthene	22.682	252	21778	0.44	ng/ul	99
25) Benzo(k)fluoranthene	22.723	252	23276	0.42	ng/ul	100
26) Benzo(a)pyrene	23.232	252	17768	0.41	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.483	276	24608	0.43	ng/ul#	82
28) Dibenzo(a,h)anthracene	25.507	278	19925	0.43	ng/ul	99
29) Benzo(g,h,i)perylene	26.147	276	20980	0.43	ng/ul	100

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

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