

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN123124\
 Data File : BN035869.D
 Acq On : 31 Dec 2024 15:15
 Operator : RC/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV243

Quant Time: Dec 31 15:51:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN123124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 31 15:06:22 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.837	152	1385	0.400	ng/ul	0.00
4) Naphthalene-d8	10.631	136	2888	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.472	164	1526	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.212	188	3142	0.400	ng/ul	0.00
17) Chrysene-d12	21.392	240	2682	0.400	ng/ul	0.00
23) Perylene-d12	23.698	264	2897	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.288	96	573	0.410	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.220	152	1508	0.402	ng/ul	0.00
18) Fluoranthene-d10	19.238	212	3014	0.425	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.322	88	618	0.406	ng/ul#	86
5) Naphthalene	10.680	128	3240	0.421	ng/ul	98
7) 2-Methylnaphthalene	12.297	142	2090	0.428	ng/ul	100
8) 1-Methylnaphthalene	12.517	142	1965	0.399	ng/ul	99
10) Acenaphthylene	14.190	152	3178	0.445	ng/ul	99
11) Acenaphthene	14.532	153	2176	0.436	ng/ul	97
12) Fluorene	15.522	166	2383	0.446	ng/ul	99
14) Pentachlorophenol	16.857	266	269	0.357	ng/ul	95
15) Phenanthrene	17.254	178	3662	0.423	ng/ul	100
16) Anthracene	17.343	178	3511	0.420	ng/ul	99
19) Fluoranthene	19.270	202	4176	0.447	ng/ul	97
20) Pyrene	19.628	202	4285	0.440	ng/ul	99
21) Benzo(a)anthracene	21.377	228	3775	0.411	ng/ul	99
22) Chrysene	21.427	228	3785	0.416	ng/ul	99
24) Benzo(b)fluoranthene	23.002	252	3669	0.421	ng/ul	99
25) Benzo(k)fluoranthene	23.049	252	3818	0.427	ng/ul	98
26) Benzo(a)pyrene	23.596	252	3452	0.421	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.058	276	4330	0.408	ng/ul#	89
28) Dibenzo(a,h)anthracene	26.078	278	3414	0.408	ng/ul	98
29) Benzo(g,h,i)perylene	26.776	276	3384	0.402	ng/ul	100

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

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