

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082621\
 Data File : BN016066.D
 Acq On : 25 Aug 2021 19:52
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV207

Quant Time: Aug 26 02:03:34 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN082621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 26 02:01:41 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.871	152	1372	0.400	ng/ul	0.00
4) Naphthalene-d8	10.666	136	5679	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.502	164	3440	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.251	188	7510	0.400	ng/ul	0.00
17) Chrysene-d12	21.436	240	7521	0.400	ng/ul	0.00
23) Perylene-d12	23.827	264	6507	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.323	96	758	0.408	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	3485	0.408	ng/ul	0.00
18) Fluoranthene-d10	19.276	212	8159	0.403	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.357	88	760	0.405	ng/ul	97
5) Naphthalene	10.716	128	6299	0.429	ng/ul	99
7) 2-Methylnaphthalene	12.333	142	4277	0.431	ng/ul	98
8) 1-Methylnaphthalene	12.547	142	3953	0.407	ng/ul	99
10) Acenaphthylene	14.224	152	5141	0.423	ng/ul	99
11) Acenaphthene	14.566	153	4424	0.422	ng/ul	99
12) Fluorene	15.552	166	5226	0.421	ng/ul	98
14) Pentachlorophenol	16.900	266	598	0.380	ng/ul	98
15) Phenanthrene	17.293	178	8658	0.418	ng/ul	99
16) Anthracene	17.381	178	7581	0.424	ng/ul	99
19) Fluoranthene	19.308	202	10015	0.426	ng/ul	99
20) Pyrene	19.671	202	10133	0.428	ng/ul	99
21) Benzo(a)anthracene	21.419	228	9858	0.421	ng/ul	99
22) Chrysene	21.474	228	10760	0.427	ng/ul	99
24) Benzo(b)fluoranthene	23.096	252	10533	0.434	ng/ul	99
25) Benzo(k)fluoranthene	23.146	252	10321	0.422	ng/ul	99
26) Benzo(a)pyrene	23.719	252	8846	0.433	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.297	276	11479	0.425	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.313	278	9654	0.425	ng/ul	99
29) Benzo(g,h,i)perylene	27.051	276	9774	0.420	ng/ul	98

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

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