

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021822\
 Data File : BN018654.D
 Acq On : 18 Feb 2022 13:55
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV208

Quant Time: Feb 18 14:51:09 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN021822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 18 13:39:55 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.917	152	6636	0.400	ng/ul	0.00
4) Naphthalene-d8	10.725	136	16474	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.551	164	7970	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.291	188	14503	0.400	ng/ul	0.00
17) Chrysene-d12	21.472	240	12339	0.400	ng/ul #	0.00
23) Perylene-d12	23.872	264	11466	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.288	96	2704	0.359	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.314	152	8462	0.376	ng/ul	0.00
18) Fluoranthene-d10	19.316	212	13696	0.372	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.326	88	3005	0.389	ng/ul#	84
5) Naphthalene	10.774	128	17753	0.381	ng/ul	99
7) 2-Methylnaphthalene	12.385	142	11459	0.381	ng/ul	100
8) 1-Methylnaphthalene	12.605	142	10651	0.355	ng/ul	100
10) Acenaphthylene	14.273	152	14202	0.395	ng/ul	100
11) Acenaphthene	14.616	153	10862	0.377	ng/ul	99
12) Fluorene	15.596	166	11536	0.373	ng/ul	98
14) Pentachlorophenol	16.945	266	1150	0.389	ng/ul	99
15) Phenanthrene	17.333	178	17731	0.372	ng/ul	99
16) Anthracene	17.422	178	15650	0.367	ng/ul	99
19) Fluoranthene	19.348	202	18801	0.363	ng/ul#	96
20) Pyrene	19.711	202	19197	0.359	ng/ul	96
21) Benzo(a)anthracene	21.458	228	17054	0.357	ng/ul	99
22) Chrysene	21.510	228	18644	0.374	ng/ul	100
24) Benzo(b)fluoranthene	23.138	252	19059	0.374	ng/ul	95
25) Benzo(k)fluoranthene	23.188	252	19041	0.374	ng/ul#	95
26) Benzo(a)pyrene	23.767	252	17976	0.400	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.365	276	21385	0.373	ng/ul#	84
28) Dibenzo(a,h)anthracene	26.379	278	17272	0.384	ng/ul#	63
29) Benzo(g,h,i)perylene	27.126	276	18412	0.375	ng/ul#	93

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

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