

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101921\
 Data File : BN017005.D
 Acq On : 19 Oct 2021 14:33
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV228

Quant Time: Oct 19 15:15:40 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN101921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 19 14:22:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.548	152	2104	0.40	ng/ul	0.00
4) Naphthalene-d8	10.321	136	8052	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.202	164	4882	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.956	188	11235	0.40	ng/ul	0.00
17) Chrysene-d12	21.167	240	12091	0.40	ng/ul	0.00
23) Perylene-d12	23.376	264	12147	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.076	96	919	0.40	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.921	152	4359	0.38	ng/ul	0.00
18) Fluoranthene-d10	18.993	212	11549	0.39	ng/ul	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.109	88	1176	0.43	ng/ul#	70
5) Naphthalene	10.370	128	8517	0.35	ng/ul	100
7) 2-Methylnaphthalene	11.998	142	5770	0.38	ng/ul	100
8) 1-Methylnaphthalene	12.218	142	5894	0.38	ng/ul	100
10) Acenaphthylene	13.915	152	7430	0.38	ng/ul	100
11) Acenaphthene	14.262	153	6607	0.39	ng/ul	99
12) Fluorene	15.256	166	7587	0.39	ng/ul	100
14) Pentachlorophenol	16.610	266	872	0.32	ng/ul	100
15) Phenanthrene	16.994	178	13680	0.38	ng/ul	99
16) Anthracene	17.087	178	11150	0.37	ng/ul	99
19) Fluoranthene	19.026	202	16072	0.39	ng/ul	97
20) Pyrene	19.388	202	16438	0.39	ng/ul	99
21) Benzo(a)anthracene	21.149	228	14606	0.37	ng/ul	100
22) Chrysene	21.202	228	17580	0.39	ng/ul	99
24) Benzo(b)fluoranthene	22.715	252	18197	0.40	ng/ul	99
25) Benzo(k)fluoranthene	22.759	252	19998	0.40	ng/ul	100
26) Benzo(a)pyrene	23.280	252	15598	0.38	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.608	276	22374	0.39	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.628	278	17786	0.39	ng/ul	99
29) Benzo(g,h,i)perylene	26.282	276	20122	0.40	ng/ul	99

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

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