

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041822\
 Data File : BN019479.D
 Acq On : 18 Apr 2022 15:20
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV229

Quant Time: Apr 18 19:11:37 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN041822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 18 15:28:08 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.818	152	4741	0.400	ng/ul	0.00
4) Naphthalene-d8	10.618	136	14181	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.457	164	8096	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.201	188	13444	0.400	ng/ul	0.00
17) Chrysene-d12	21.395	240	8232	0.400	ng/ul	0.00
23) Perylene-d12	23.742	264	5740	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.173	96	2335	0.406	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.219	152	8206	0.431	ng/ul	0.00
18) Fluoranthene-d10	19.231	212	15030	0.448	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.211	88	2150	0.366	ng/ul#	62
5) Naphthalene	10.668	128	17563	0.436	ng/ul	99
7) 2-Methylnaphthalene	12.290	142	10889	0.434	ng/ul#	100
8) 1-Methylnaphthalene	12.505	142	10976	0.406	ng/ul#	100
10) Acenaphthylene	14.179	152	14163	0.475	ng/ul	99
11) Acenaphthene	14.522	153	12629	0.430	ng/ul	99
12) Fluorene	15.507	166	13002	0.431	ng/ul	99
14) Pentachlorophenol	16.872	266	622	0.347	ng/ul	87
15) Phenanthrene	17.244	178	19172	0.419	ng/ul	99
16) Anthracene	17.341	178	14015	0.426	ng/ul	99
19) Fluoranthene	19.263	202	19652	0.430	ng/ul	99
20) Pyrene	19.626	202	20978	0.424	ng/ul	99
21) Benzo(a)anthracene	21.380	228	9702	0.412	ng/ul	100
22) Chrysene	21.430	228	13538	0.442	ng/ul	99
24) Benzo(b)fluoranthene	23.029	252	10978	0.403	ng/ul	94
25) Benzo(k)fluoranthene	23.076	252	12581	0.423	ng/ul	99
26) Benzo(a)pyrene	23.640	252	10187	0.431	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	26.172	276	10868	0.451	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.199	278	7887	0.493	ng/ul#	89
29) Benzo(g,h,i)perylene	26.910	276	11302	0.443	ng/ul	99

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

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