

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112622\
 Data File : BN022880.D
 Acq On : 26 Nov 2022 13:36
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV215

Quant Time: Nov 28 03:20:28 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN112622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 28 03:16:56 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.069	152	7250	0.400	ng/u1	0.00
4) Naphthalene-d8	10.892	136	19631	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.710	164	10810	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.453	188	23307	0.400	ng/u1	0.00
17) Chrysene-d12	21.638	240	14933	0.400	ng/u1	0.00
23) Perylene-d12	24.126	264	9463	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.412	96	3539	0.407	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.476	152	11390	0.381	ng/u1	0.00
18) Fluoranthene-d10	19.480	212	22075	0.400	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.450	88	3657	0.403	ng/u1	96
5) Naphthalene	10.942	128	24676	0.414	ng/u1	100
7) 2-Methylnaphthalene	12.547	142	16168	0.434	ng/u1	100
8) 1-Methylnaphthalene	12.767	142	15521	0.410	ng/u1	99
10) Acenaphthylene	14.432	152	21851	0.410	ng/u1	100
11) Acenaphthene	14.775	153	17388	0.412	ng/u1	99
12) Fluorene	15.755	166	19858	0.416	ng/u1	100
14) Pentachlorophenol	17.099	266	1609	0.330	ng/u1#	100
15) Phenanthrene	17.496	178	32695	0.409	ng/u1	100
16) Anthracene	17.588	178	27656	0.403	ng/u1	100
19) Fluoranthene	19.508	202	32877	0.426	ng/u1	99
20) Pyrene	19.871	202	31709	0.420	ng/u1	99
21) Benzo(a)anthracene	21.620	228	22623	0.396	ng/u1	99
22) Chrysene	21.673	228	25730	0.411	ng/u1	99
24) Benzo(b)fluoranthene	23.362	252	20179	0.421	ng/u1	99
25) Benzo(k)fluoranthene	23.415	252	21215	0.419	ng/u1	99
26) Benzo(a)pyrene	24.012	252	14613	0.400	ng/u1	99
27) Indeno(1,2,3-cd)pyrene	26.740	276	16061	0.370	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.766	278	13391	0.388	ng/u1	99
29) Benzo(g,h,i)perylene	27.541	276	15157	0.400	ng/u1	100

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

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