

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042023\
 Data File : BN025027.D
 Acq On : 20 Apr 2023 07:45
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV215

Quant Time: Apr 20 08:18:19 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN042023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 20 07:57:39 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.930	152	6992	0.400	ng/u1	0.00
4) Naphthalene-d8	10.743	136	21564	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.575	164	12173	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.321	188	26677	0.400	ng/u1	0.00
17) Chrysene-d12	21.511	240	21717	0.400	ng/u1	0.00
23) Perylene-d12	23.955	264	18401	0.400	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.310	96	2588	0.357	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.332	152	11810	0.427	ng/u1	0.00
18) Fluoranthene-d10	19.349	212	26779	0.468	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.343	88	2693	0.345	ng/u1	94
5) Naphthalene	10.792	128	21871	0.396	ng/u1	100
7) 2-Methylnaphthalene	12.403	142	14242	0.421	ng/u1	100
8) 1-Methylnaphthalene	12.623	142	13547	0.376	ng/u1	100
10) Acenaphthylene	14.297	152	19536	0.431	ng/u1	100
11) Acenaphthene	14.635	153	15650	0.409	ng/u1	99
12) Fluorene	15.625	166	17693	0.414	ng/u1	99
14) Pentachlorophenol	16.980	266	1688	0.322	ng/u1	99
15) Phenanthrene	17.364	178	29686	0.394	ng/u1	100
16) Anthracene	17.457	178	24460	0.397	ng/u1	100
19) Fluoranthene	19.382	202	33321	0.424	ng/u1	99
20) Pyrene	19.744	202	34402	0.426	ng/u1	99
21) Benzo(a)anthracene	21.497	228	26880	0.411	ng/u1	99
22) Chrysene	21.546	228	30279	0.424	ng/u1	100
24) Benzo(b)fluoranthene	23.204	252	29923	0.401	ng/u1	97
25) Benzo(k)fluoranthene	23.253	252	29519	0.411	ng/u1	100
26) Benzo(a)pyrene	23.853	252	25211	0.408	ng/u1#	95
27) Indeno(1,2,3-cd)pyrene	26.507	276	28746	0.404	ng/u1#	99
28) Dibenzo(a,h)anthracene	26.527	278	22358	0.402	ng/u1	99
29) Benzo(g,h,i)perylene	27.285	276	23751	0.392	ng/u1	99

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

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