

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN051624\
 Data File : BN031535.D
 Acq On : 15 May 2024 13:06
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV222

Quant Time: May 15 13:35:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN051624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 15 12:57:23 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.700	152	10751	0.400	ng/u1	0.00
4) Naphthalene-d8	10.481	136	33261	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.336	164	17061	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.079	188	37560	0.400	ng/u1	0.00
17) Chrysene-d12	21.264	240	30638	0.400	ng/u1	0.00
23) Perylene-d12	23.500	264	27352	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.207	96	5342	0.412	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.076	152	18304	0.389	ng/u1	0.00
18) Fluoranthene-d10	19.106	212	34451	0.405	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.240	88	5663	0.420	ng/u1	94
5) Naphthalene	10.530	128	36469	0.410	ng/u1	100
7) 2-Methylnaphthalene	12.147	142	23154	0.412	ng/u1	100
8) 1-Methylnaphthalene	12.367	142	21592	0.364	ng/u1	100
10) Acenaphthylene	14.054	152	31419	0.406	ng/u1	100
11) Acenaphthene	14.397	153	23342	0.396	ng/u1	99
12) Fluorene	15.387	166	26568	0.406	ng/u1	99
14) Pentachlorophenol	16.721	266	2027	0.346	ng/u1	98
15) Phenanthrene	17.122	178	43911	0.399	ng/u1	100
16) Anthracene	17.210	178	38682	0.417	ng/u1	99
19) Fluoranthene	19.138	202	46734	0.410	ng/u1	100
20) Pyrene	19.501	202	49219	0.437	ng/u1	100
21) Benzo(a)anthracene	21.250	228	38806	0.366	ng/u1	100
22) Chrysene	21.299	228	46376	0.405	ng/u1	99
24) Benzo(b)fluoranthene	22.828	252	39479	0.392	ng/u1	98
25) Benzo(k)fluoranthene	22.874	252	42274	0.388	ng/u1	98
26) Benzo(a)pyrene	23.403	252	36394	0.412	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	25.764	276	40126	0.391	ng/u1	100
28) Dibenzo(a,h)anthracene	25.780	278	31333	0.385	ng/u1	99
29) Benzo(g,h,i)perylene	26.451	276	31315	0.380	ng/u1	99

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

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