

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN012125\
 Data File : BN036008.D
 Acq On : 21 Jan 2025 16:41
 Operator : RC/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV201

Quant Time: Jan 21 23:53:09 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN012125.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 21 23:52:22 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.812	152	2858	0.400	ng/ul	0.00
4) Naphthalene-d8	10.605	136	5645	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.451	164	2812	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.192	188	5615	0.400	ng/ul	0.00
17) Chrysene-d12	21.374	240	4425	0.400	ng/ul	0.00
23) Perylene-d12	23.674	264	4716	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	1277	0.408	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.194	152	2833	0.397	ng/ul	0.00
18) Fluoranthene-d10	19.216	212	4906	0.403	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.310	88	1338	0.401	ng/ul	95
5) Naphthalene	10.655	128	6259	0.402	ng/ul	99
7) 2-Methylnaphthalene	12.271	142	3880	0.402	ng/ul	100
8) 1-Methylnaphthalene	12.491	142	3612	0.368	ng/ul	99
10) Acenaphthylene	14.169	152	5278	0.398	ng/ul	98
11) Acenaphthene	14.511	153	3887	0.398	ng/ul	99
12) Fluorene	15.501	166	4146	0.385	ng/ul	98
14) Pentachlorophenol	16.842	266	786	0.400	ng/ul	99
15) Phenanthrene	17.234	178	6535	0.397	ng/ul	99
16) Anthracene	17.323	178	6034	0.395	ng/ul	99
19) Fluoranthene	19.249	202	6734	0.400	ng/ul	100
20) Pyrene	19.611	202	6975	0.397	ng/ul	99
21) Benzo(a)anthracene	21.359	228	6207	0.393	ng/ul	99
22) Chrysene	21.409	228	6425	0.398	ng/ul	99
24) Benzo(b)fluoranthene	22.978	252	6352	0.404	ng/ul	99
25) Benzo(k)fluoranthene	23.025	252	6588	0.398	ng/ul	99
26) Benzo(a)pyrene	23.569	252	5919	0.416	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.010	276	7521	0.400	ng/ul	100
28) Dibenzo(a,h)anthracene	26.031	278	5936	0.410	ng/ul	99
29) Benzo(g,h,i)perylene	26.732	276	5725	0.392	ng/ul	98

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

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