

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082724\
 Data File : BN033636.D
 Acq On : 27 Aug 2024 14:48
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV208

Quant Time: Aug 27 23:49:26 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN082724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 23:48:09 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.540	152	5585	0.400	ng/ul	0.00
4) Naphthalene-d8	10.299	136	15133	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.179	164	7741	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.927	188	15980	0.400	ng/ul	0.00
17) Chrysene-d12	21.138	240	13187	0.400	ng/ul	0.00
23) Perylene-d12	23.301	264	14416	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.148	96	2508	0.437	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.900	152	8426	0.388	ng/ul	0.00
18) Fluoranthene-d10	18.966	212	15481	0.394	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.177	88	2572	0.404	ng/ul	96
5) Naphthalene	10.349	128	16358	0.404	ng/ul	100
7) 2-Methylnaphthalene	11.971	142	10482	0.409	ng/ul	100
8) 1-Methylnaphthalene	12.197	142	9720	0.369	ng/ul	99
10) Acenaphthylene	13.892	152	14575	0.415	ng/ul	99
11) Acenaphthene	14.239	153	10391	0.409	ng/ul	100
12) Fluorene	15.234	166	11683	0.412	ng/ul	99
14) Pentachlorophenol	16.585	266	1645	0.327	ng/ul	99
15) Phenanthrene	16.969	178	18217	0.386	ng/ul	99
16) Anthracene	17.058	178	16833	0.397	ng/ul	100
19) Fluoranthene	18.994	202	20341	0.391	ng/ul	98
20) Pyrene	19.361	202	21458	0.388	ng/ul	99
21) Benzo(a)anthracene	21.124	228	19890	0.383	ng/ul	99
22) Chrysene	21.173	228	20020	0.382	ng/ul	99
24) Benzo(b)fluoranthene	22.661	252	21189	0.379	ng/ul	99
25) Benzo(k)fluoranthene	22.702	252	20910	0.374	ng/ul	99
26) Benzo(a)pyrene	23.207	252	19138	0.426	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.452	276	25381	0.376	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.468	278	19908	0.380	ng/ul	99
29) Benzo(g,h,i)perylene	26.102	276	20033	0.380	ng/ul	99

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

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