

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091624\
 Data File : BN033986.D
 Acq On : 16 Sep 2024 13:55
 Operator : JU/RC
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4310

Quant Time: Sep 16 14:29:47 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 : Sun Sep 15 01:52:43 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.477	152	6683	0.400	ng/ul	0.00
4) Naphthalene-d8	10.233	136	18228	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.123	164	9151	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.877	188	18517	0.400	ng/ul	0.00
17) Chrysene-d12	21.095	240	15718	0.400	ng/ul	0.00
23) Perylene-d12	23.234	264	17822	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.097	96	2992	0.435	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.834	152	10568	0.404	ng/ul	-0.01
18) Fluoranthene-d10	18.920	212	19562	0.417	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.131	88	3324	0.436	ng/ul	93
5) Naphthalene	10.283	128	19665	0.403	ng/ul	100
7) 2-Methylnaphthalene	11.911	142	12336	0.399	ng/ul	100
8) 1-Methylnaphthalene	12.131	142	12775	0.402	ng/ul	99
10) Acenaphthylene	13.837	152	16636	0.400	ng/ul	100
11) Acenaphthene	14.184	153	12399	0.413	ng/ul	99
12) Fluorene	15.178	166	13885	0.414	ng/ul	99
14) Pentachlorophenol	16.535	266	1950	0.335	ng/ul	99
15) Phenanthrene	16.919	178	21341	0.391	ng/ul	99
16) Anthracene	17.007	178	19334	0.394	ng/ul	99
19) Fluoranthene	18.948	202	24314	0.392	ng/ul	98
20) Pyrene	19.315	202	25750	0.391	ng/ul	100
21) Benzo(a)anthracene	21.077	228	24511	0.396	ng/ul	100
22) Chrysene	21.130	228	25075	0.401	ng/ul	99
24) Benzo(b)fluoranthene	22.602	252	25853	0.374	ng/ul	97
25) Benzo(k)fluoranthene	22.643	252	27181	0.393	ng/ul	98
26) Benzo(a)pyrene	23.140	252	21910	0.394	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.344	276	32284	0.387	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.364	278	25116	0.388	ng/ul	98
29) Benzo(g,h,i)perylene	25.992	276	25566	0.392	ng/ul	99

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

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