

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034977.D
 Acq On : 10 Nov 2024 13:56
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 71 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4329

Quant Time: Nov 10 21:52:21 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 09 00:09:12 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.566	152	2506	0.400	ng/ul	0.00
4) Naphthalene-d8	10.328	136	5947	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.199	164	4323	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.950	188	10034	0.400	ng/ul	0.00
17) Chrysene-d12	21.142	240	9539	0.400	ng/ul	0.00
23) Perylene-d12	23.305	264	10188	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	944	0.396	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.928	152	3579	0.393	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	10100	0.412	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.182	88	1019	0.390	ng/ul	96
5) Naphthalene	10.377	128	5934	0.399	ng/ul	98
7) 2-Methylnaphthalene	12.000	142	4263	0.384	ng/ul	99
8) 1-Methylnaphthalene	12.220	142	4328	0.388	ng/ul	98
10) Acenaphthylene	13.916	152	6714	0.380	ng/ul	99
11) Acenaphthene	14.263	153	5184	0.387	ng/ul	100
12) Fluorene	15.253	166	6228	0.366	ng/ul	99
14) Pentachlorophenol	16.603	266	786	0.339	ng/ul	98
15) Phenanthrene	16.987	178	10552	0.394	ng/ul	99
16) Anthracene	17.080	178	9711	0.385	ng/ul	99
19) Fluoranthene	19.009	202	12896	0.405	ng/ul#	97
20) Pyrene	19.372	202	13173	0.399	ng/ul#	96
21) Benzo(a)anthracene	21.125	228	12945	0.375	ng/ul	100
22) Chrysene	21.177	228	13487	0.383	ng/ul	99
24) Benzo(b)fluoranthene	22.662	252	13373	0.373	ng/ul	99
25) Benzo(k)fluoranthene	22.703	252	14137	0.377	ng/ul	99
26) Benzo(a)pyrene	23.209	252	12190	0.381	ng/ul#	97
27) Indeno(1,2,3-cd)pyrene	25.456	276	16190	0.414	ng/ul#	95
28) Dibenzo(a,h)anthracene	25.473	278	12671	0.416	ng/ul	98
29) Benzo(g,h,i)perylene	26.111	276	12617	0.422	ng/ul	98

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

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