

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121024\
 Data File : BN035577.D
 Acq On : 11 Dec 2024 22:06
 Operator : RC/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4362

Quant Time: Dec 11 22:33:14 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 10 16:10:57 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.288	152	2486	0.400	ng/ul	0.00
4) Naphthalene-d8	10.036	136	6921	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.949	164	4669	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.717	188	10005	0.400	ng/ul	0.00
17) Chrysene-d12	20.961	240	8121	0.400	ng/ul	0.00
23) Perylene-d12	23.054	264	8317	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.963	96	973	0.428	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.637	152	4142	0.408	ng/ul	0.00
18) Fluoranthene-d10	18.768	212	9309	0.416	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.993	88	1045	0.424	ng/ul#	91
5) Naphthalene	10.086	128	7014	0.402	ng/ul	99
7) 2-Methylnaphthalene	11.714	142	4871	0.403	ng/ul	98
8) 1-Methylnaphthalene	11.939	142	5008	0.410	ng/ul	98
10) Acenaphthylene	13.662	152	7642	0.413	ng/ul	100
11) Acenaphthene	14.014	153	5903	0.417	ng/ul	99
12) Fluorene	15.017	166	6672	0.401	ng/ul	99
14) Pentachlorophenol	16.380	266	951	0.408	ng/ul	98
15) Phenanthrene	16.759	178	10880	0.411	ng/ul	100
16) Anthracene	16.848	178	9582	0.390	ng/ul	100
19) Fluoranthene	18.800	202	11832	0.418	ng/ul	98
20) Pyrene	19.167	202	12465	0.424	ng/ul#	96
21) Benzo(a)anthracene	20.947	228	11038	0.392	ng/ul	100
22) Chrysene	20.996	228	11799	0.416	ng/ul	99
24) Benzo(b)fluoranthene	22.440	252	10983	0.400	ng/ul	97
25) Benzo(k)fluoranthene	22.481	252	12278	0.433	ng/ul	97
26) Benzo(a)pyrene	22.963	252	10041	0.400	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.084	276	12494	0.368	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.101	278	8829	0.332	ng/ul	97
29) Benzo(g,h,i)perylene	25.698	276	8967	0.335	ng/ul	98

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

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