

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112624\
 Data File : BN035318.D
 Acq On : 26 Nov 2024 17:07
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4353

Quant Time: Nov 26 22:14:08 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN112624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 26 06:09:15 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.314	152	3358	0.400	ng/ul	0.00
4) Naphthalene-d8	10.058	136	8647	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.972	164	4975	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.734	188	9193	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.979	240	5730	0.400	ng/ul	# 0.00
23) Perylene-d12	23.074	264	6161	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.976	96	1358	0.442	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.659	152	4660	0.368	ng/ul	0.00
18) Fluoranthene-d10	18.786	212	7315	0.463	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.005	88	1388	0.417	ng/ul	93
5) Naphthalene	10.108	128	8891	0.408	ng/ul	99
7) 2-Methylnaphthalene	11.736	142	5902	0.391	ng/ul	99
8) 1-Methylnaphthalene	11.961	142	5987	0.393	ng/ul	99
10) Acenaphthylene	13.680	152	8445	0.428	ng/ul	99
11) Acenaphthene	14.032	153	6395	0.424	ng/ul	97
12) Fluorene	15.036	166	6962	0.393	ng/ul	98
14) Pentachlorophenol	16.396	266	660	0.308	ng/ul	97
15) Phenanthrene	16.776	178	10225	0.421	ng/ul	99
16) Anthracene	16.869	178	8869	0.393	ng/ul	99
19) Fluoranthene	18.819	202	9774	0.490	ng/ul#	95
20) Pyrene	19.186	202	9706	0.468	ng/ul#	93
21) Benzo(a)anthracene	20.961	228	7821	0.393	ng/ul	99
22) Chrysene	21.014	228	8591	0.429	ng/ul	100
24) Benzo(b)fluoranthene	22.457	252	8409	0.413	ng/ul	96
25) Benzo(k)fluoranthene	22.498	252	9227	0.439	ng/ul#	96
26) Benzo(a)pyrene	22.983	252	7729	0.415	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	25.114	276	10270	0.409	ng/ul#	87
28) Dibenzo(a,h)anthracene	25.131	278	7751	0.393	ng/ul#	96
29) Benzo(g,h,i)perylene	25.728	276	8464	0.427	ng/ul#	95

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

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