

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112624\
 Data File : BN035306.D
 Acq On : 26 Nov 2024 08:08
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4352

Quant Time: Nov 26 08:31:19 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 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	2932	0.400	ng/ul	0.00
4) Naphthalene-d8	10.058	136	6911	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.972	164	3728	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.738	188	7604	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.979	240	6417	0.400	ng/ul	# 0.00
23) Perylene-d12	23.074	264	7260	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.976	96	1162	0.434	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.664	152	3536	0.349	ng/ul	0.00
18) Fluoranthene-d10	18.791	212	7118	0.402	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.010	88	1297	0.446	ng/ul	97
5) Naphthalene	10.108	128	7028	0.404	ng/ul	99
7) 2-Methylnaphthalene	11.741	142	4532	0.375	ng/ul	99
8) 1-Methylnaphthalene	11.967	142	4576	0.375	ng/ul	99
10) Acenaphthylene	13.685	152	6406	0.433	ng/ul	99
11) Acenaphthene	14.037	153	4889	0.432	ng/ul	97
12) Fluorene	15.036	166	5364	0.404	ng/ul	98
14) Pentachlorophenol	16.396	266	593	0.335	ng/ul	99
15) Phenanthrene	16.781	178	8588	0.427	ng/ul	99
16) Anthracene	16.869	178	7425	0.398	ng/ul	99
19) Fluoranthene	18.819	202	9591	0.429	ng/ul#	94
20) Pyrene	19.186	202	9880	0.425	ng/ul#	91
21) Benzo(a)anthracene	20.964	228	8783	0.394	ng/ul	99
22) Chrysene	21.014	228	9801	0.437	ng/ul	99
24) Benzo(b)fluoranthene	22.460	252	9344	0.389	ng/ul	95
25) Benzo(k)fluoranthene	22.501	252	10781	0.436	ng/ul#	95
26) Benzo(a)pyrene	22.983	252	8724	0.398	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	25.114	276	12031	0.406	ng/ul#	85
28) Dibenzo(a,h)anthracene	25.135	278	9083	0.391	ng/ul#	94
29) Benzo(g,h,i)perylene	25.732	276	9252	0.396	ng/ul#	94

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

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