

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121324\
 Data File : BN035599.D
 Acq On : 13 Dec 2024 16:32
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4366

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/14/2024
 Supervised By :mohammad ahmed 12/14/2024

Quant Time: Dec 13 17:17:05 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	7.297	152	1651	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.042	136	5135	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	13.958	164	3570	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.726	188	7962	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.970	240	6910	0.400	ng/ul	# 0.00
23) Perylene-d12	23.062	264	7819	0.400	ng/ul	# 0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.963	96	594m	0.394	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.648	152	3046	0.405	ng/ul	0.00
18) Fluoranthene-d10	18.777	212	8396	0.440	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.993	88	685	0.418	ng/ul#	9
5) Naphthalene	10.091	128	4932m	0.381	ng/ul	
7) 2-Methylnaphthalene	11.719	142	3524	0.393	ng/ul	97
8) 1-Methylnaphthalene	11.950	142	3635	0.401	ng/ul	98
10) Acenaphthylene	13.671	152	5841	0.412	ng/ul	98
11) Acenaphthene	14.018	153	4368	0.403	ng/ul	96
12) Fluorene	15.022	166	5044	0.397	ng/ul	99
14) Pentachlorophenol	16.392	266	1028	0.554	ng/ul	96
15) Phenanthrene	16.768	178	8246	0.392	ng/ul	97
16) Anthracene	16.857	178	8135	0.416	ng/ul	97
19) Fluoranthene	18.809	202	10083	0.419	ng/ul#	96
20) Pyrene	19.176	202	11722	0.468	ng/ul#	93
21) Benzo(a)anthracene	20.955	228	9692	0.404	ng/ul#	85
22) Chrysene	21.005	228	9594	0.398	ng/ul#	85
24) Benzo(b)fluoranthene	22.448	252	9762	0.378	ng/ul#	44
25) Benzo(k)fluoranthene	22.489	252	10507	0.394	ng/ul#	37
26) Benzo(a)pyrene	22.972	252	9356	0.396	ng/ul#	30
27) Indeno(1,2,3-cd)pyrene	25.104	276	11109m	0.348	ng/ul	
28) Dibenzo(a,h)anthracene	25.121	278	9021m	0.361	ng/ul	
29) Benzo(g,h,i)perylene	25.718	276	8899	0.354	ng/ul#	43

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

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