

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111924\
 Data File : BN035189.D
 Acq On : 20 Nov 2024 08:04
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4345

Quant Time: Nov 20 09:38:18 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 : Mon Nov 18 00:44:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.541	152	3680	0.400	ng/ul	0.00
4) Naphthalene-d8	10.306	136	9095	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.185	164	6214	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.937	188	13002	0.400	ng/ul	0.00
17) Chrysene-d12	21.145	240	9390	0.400	ng/ul	0.00
23) Perylene-d12	23.317	264	9800	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.136	96	1385	0.412	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.901	152	5573	0.418	ng/ul	-0.01
18) Fluoranthene-d10	18.977	212	11751	0.454	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.170	88	1525	0.418	ng/ul	97
5) Naphthalene	10.355	128	9415	0.411	ng/ul	99
7) 2-Methylnaphthalene	11.978	142	6721	0.423	ng/ul	100
8) 1-Methylnaphthalene	12.198	142	6754	0.421	ng/ul	99
10) Acenaphthylene	13.898	152	9921	0.402	ng/ul	100
11) Acenaphthene	14.245	153	7716	0.409	ng/ul	99
12) Fluorene	15.240	166	8903	0.402	ng/ul	99
14) Pentachlorophenol	16.595	266	930	0.307	ng/ul	97
15) Phenanthrene	16.979	178	13890	0.404	ng/ul	99
16) Anthracene	17.068	178	12613	0.395	ng/ul	100
19) Fluoranthene	19.005	202	14893	0.456	ng/ul	99
20) Pyrene	19.372	202	15166	0.446	ng/ul	99
21) Benzo(a)anthracene	21.131	228	13024	0.400	ng/ul	100
22) Chrysene	21.180	228	13461	0.411	ng/ul	99
24) Benzo(b)fluoranthene	22.671	252	13040	0.403	ng/ul	98
25) Benzo(k)fluoranthene	22.712	252	14268	0.427	ng/ul	98
26) Benzo(a)pyrene	23.220	252	12080	0.408	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.467	276	16527	0.413	ng/ul	99
28) Dibenzo(a,h)anthracene	25.487	278	12949	0.413	ng/ul	98
29) Benzo(g,h,i)perylene	26.124	276	12866	0.408	ng/ul	99

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

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