

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111924\
 Data File : BN035164.D
 Acq On : 19 Nov 2024 16:28
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4343

Quant Time: Nov 19 17:13:48 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	7.545	152	3035	0.400	ng/ul	0.00
4) Naphthalene-d8	10.306	136	7523	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.185	164	5577	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.941	188	13051	0.400	ng/ul	0.00
17) Chrysene-d12	21.148	240	9993	0.400	ng/ul	0.00
23) Perylene-d12	23.320	264	9110	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	1113	0.401	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.906	152	4761	0.432	ng/ul	0.00
18) Fluoranthene-d10	18.977	212	12764	0.463	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.170	88	1262	0.419	ng/ul	97
5) Naphthalene	10.355	128	7855	0.414	ng/ul	99
7) 2-Methylnaphthalene	11.978	142	5655	0.430	ng/ul	99
8) 1-Methylnaphthalene	12.203	142	5779	0.436	ng/ul	98
10) Acenaphthylene	13.902	152	8900	0.402	ng/ul	99
11) Acenaphthene	14.249	153	6945	0.410	ng/ul	99
12) Fluorene	15.244	166	8385	0.422	ng/ul	99
14) Pentachlorophenol	16.595	266	1020	0.336	ng/ul	98
15) Phenanthrene	16.979	178	14002	0.406	ng/ul	99
16) Anthracene	17.072	178	12727	0.397	ng/ul	99
19) Fluoranthene	19.009	202	16173	0.465	ng/ul	99
20) Pyrene	19.372	202	16513	0.456	ng/ul	99
21) Benzo(a)anthracene	21.131	228	13861	0.400	ng/ul	100
22) Chrysene	21.183	228	14239	0.408	ng/ul	99
24) Benzo(b)fluoranthene	22.671	252	12634	0.420	ng/ul	99
25) Benzo(k)fluoranthene	22.714	252	13323	0.429	ng/ul	99
26) Benzo(a)pyrene	23.223	252	11020	0.400	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.473	276	13979	0.376	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.490	278	10928	0.375	ng/ul	99
29) Benzo(g,h,i)perylene	26.127	276	10769	0.367	ng/ul	99

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

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