

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034961.D
 Acq On : 10 Nov 2024 03:48
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4328

Quant Time: Nov 10 04:11:07 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 09 00:09:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.566	152	3265	0.400	ng/ul	0.00
4) Naphthalene-d8	10.333	136	8143	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.203	164	6354	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.949	188	15806	0.400	ng/ul	0.00
17) Chrysene-d12	21.142	240	14876	0.400	ng/ul	0.00
23) Perylene-d12	23.305	264	13667	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	1150	0.370	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.928	152	5103	0.409	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	16456	0.431	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.187	88	1249	0.367	ng/ul#	94
5) Naphthalene	10.383	128	8119	0.398	ng/ul	98
7) 2-Methylnaphthalene	12.005	142	6067	0.399	ng/ul	100
8) 1-Methylnaphthalene	12.225	142	6136	0.402	ng/ul	99
10) Acenaphthylene	13.921	152	9986	0.385	ng/ul	99
11) Acenaphthene	14.263	153	7705	0.392	ng/ul	98
12) Fluorene	15.258	166	9428	0.377	ng/ul	100
14) Pentachlorophenol	16.603	266	1169	0.320	ng/ul	97
15) Phenanthrene	16.992	178	16563	0.392	ng/ul	98
16) Anthracene	17.080	178	15309	0.386	ng/ul	98
19) Fluoranthene	19.014	202	20938	0.421	ng/ul	99
20) Pyrene	19.376	202	21683	0.421	ng/ul	99
21) Benzo(a)anthracene	21.125	228	20370	0.378	ng/ul	99
22) Chrysene	21.177	228	20728	0.377	ng/ul	99
24) Benzo(b)fluoranthene	22.662	252	18554	0.386	ng/ul	99
25) Benzo(k)fluoranthene	22.703	252	19761	0.393	ng/ul	98
26) Benzo(a)pyrene	23.211	252	16208	0.378	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.456	276	19798	0.377	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.473	278	15527	0.380	ng/ul	98
29) Benzo(g,h,i)perylene	26.114	276	15190	0.379	ng/ul	99

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

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