

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034916.D
 Acq On : 08 Nov 2024 22:28
 Operator : RC/JU
 Sample : SSTD0.834
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.8219

Quant Time: Nov 08 23:56:26 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 : Fri Nov 08 23:54:11 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	1702	0.400	ng/u1	0.00
4) Naphthalene-d8	10.339	136	4088	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.208	164	3078	0.400	ng/u1	0.00
13) Phenanthrene-d10	16.954	188	7911	0.400	ng/u1	0.00
17) Chrysene-d12	21.148	240	8518	0.400	ng/u1	0.00
23) Perylene-d12	23.314	264	8682	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	1290	0.732	ng/u1	0.00
6) 2-Methylnaphthalene-d10	11.934	152	4970	0.970	ng/u1	0.00
18) Fluoranthene-d10	18.986	212	17116	0.617	ng/u1	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.182	88	1420	0.734	ng/u1	94
5) Naphthalene	10.388	128	8099	0.748	ng/u1	98
7) 2-Methylnaphthalene	12.011	142	6047	0.859	ng/u1	99
8) 1-Methylnaphthalene	12.231	142	6088	0.849	ng/u1	98
10) Acenaphthylene	13.926	152	9876	0.705	ng/u1	98
11) Acenaphthene	14.273	153	7547	0.704	ng/u1	99
12) Fluorene	15.263	166	9578	0.802	ng/u1	100
14) Pentachlorophenol	16.608	266	1467	0.940	ng/u1	97
15) Phenanthrene	16.996	178	16739	0.700	ng/u1	98
16) Anthracene	17.085	178	15826	0.715	ng/u1	98
19) Fluoranthene	19.019	202	22518	0.545	ng/u1	100
20) Pyrene	19.381	202	23168	0.539	ng/u1	100
21) Benzo(a)anthracene	21.131	228	24301	0.713	ng/u1	99
22) Chrysene	21.183	228	24658	0.722	ng/u1	100
24) Benzo(b)fluoranthene	22.671	252	24310	0.688	ng/u1	97
25) Benzo(k)fluoranthene	22.712	252	25388	0.712	ng/u1	96
26) Benzo(a)pyrene	23.220	252	21571	0.709	ng/u1	95
27) Indeno(1,2,3-cd)pyrene	25.467	276	26516	0.818	ng/u1	99
28) Dibenzo(a,h)anthracene	25.487	278	20809	0.850	ng/u1	97
29) Benzo(g,h,i)perylene	26.124	276	20245	0.795	ng/u1	99

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

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