

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062724\
 Data File : BN032275.D
 Acq On : 27 Jun 2024 11:06
 Operator : MA/JU
 Sample : SSTD0.403
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4232

Quant Time: Jun 27 12:50:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN062724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 27 12:48:23 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.620	152	7662	0.400	ng/u1	0.00
4) Naphthalene-d8	10.398	136	23860	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.267	164	15976	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.025	188	37608	0.400	ng/u1	0.00
17) Chrysene-d12	21.226	240	31365	0.400	ng/u1	0.00
23) Perylene-d12	23.453	264	31845	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.156	96	3295	0.405	ng/u1	0.00
6) 2-Methylnaphthalene-d10	11.999	152	13178	0.352	ng/u1	0.00
18) Fluoranthene-d10	19.059	212	34780	0.373	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.186	88	3740	0.410	ng/u1	100
5) Naphthalene	10.448	128	23383	0.377	ng/u1	100
7) 2-Methylnaphthalene	12.076	142	15446	0.361	ng/u1	100
8) 1-Methylnaphthalene	12.296	142	16046	0.362	ng/u1	100
10) Acenaphthylene	13.989	152	23992	0.362	ng/u1	100
11) Acenaphthene	14.332	153	18720	0.375	ng/u1	100
12) Fluorene	15.326	166	22313	0.368	ng/u1	100
14) Pentachlorophenol	16.699	266	2828	0.327	ng/u1	100
15) Phenanthrene	17.067	178	37196	0.361	ng/u1	100
16) Anthracene	17.160	178	32686	0.354	ng/u1	100
19) Fluoranthene	19.092	202	44119	0.386	ng/u1	100
20) Pyrene	19.454	202	45983	0.384	ng/u1	100
21) Benzo(a)anthracene	21.214	228	39819	0.342	ng/u1	100
22) Chrysene	21.267	228	44003	0.373	ng/u1	100
24) Benzo(b)fluoranthene	22.784	252	45721	0.336	ng/u1	100
25) Benzo(k)fluoranthene	22.828	252	45329	0.345	ng/u1	100
26) Benzo(a)pyrene	23.360	252	33908	0.351	ng/u1	100
27) Indeno(1,2,3-cd)pyrene	25.700	276	50606	0.379	ng/u1	100
28) Dibenzo(a,h)anthracene	25.710	278	38068	0.371	ng/u1	100
29) Benzo(g,h,i)perylene	26.381	276	40235	0.379	ng/u1	100

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

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