

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101921\
 Data File : BN017001.D
 Acq On : 19 Oct 2021 12:01
 Operator : CG/JU
 Sample : SST0.424
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.4224

Quant Time: Oct 19 12:51:53 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN101921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 19 12:51:41 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.548	152	2193	0.40	ng/ul	0.00
4) Naphthalene-d8	10.321	136	8319	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.197	164	4974	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.952	188	11210	0.40	ng/ul	0.00
17) Chrysene-d12	21.164	240	12439	0.40	ng/ul	0.00
23) Perylene-d12	23.376	264	12913	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.076	96	950	0.37	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.921	152	4499	0.38	ng/ul	0.00
18) Fluoranthene-d10	18.993	212	11932	0.36	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.109	88	1238	0.44	ng/ul	100
5) Naphthalene	10.370	128	9618	0.43	ng/ul	100
7) 2-Methylnaphthalene	11.998	142	5942	0.41	ng/ul	100
8) 1-Methylnaphthalene	12.218	142	6083	0.42	ng/ul	100
10) Acenaphthylene	13.915	152	7717	0.41	ng/ul	100
11) Acenaphthene	14.262	153	6674	0.42	ng/ul	100
12) Fluorene	15.257	166	7720	0.43	ng/ul	100
14) Pentachlorophenol	16.610	266	1035	0.57	ng/ul	100
15) Phenanthrene	16.994	178	13737	0.43	ng/ul	100
16) Anthracene	17.087	178	11512	0.41	ng/ul	100
19) Fluoranthene	19.021	202	16330	0.39	ng/ul	100
20) Pyrene	19.388	202	16846	0.40	ng/ul	100
21) Benzo(a)anthracene	21.146	228	15481	0.41	ng/ul	100
22) Chrysene	21.199	228	17949	0.42	ng/ul	100
24) Benzo(b)fluoranthene	22.710	252	19012	0.39	ng/ul	100
25) Benzo(k)fluoranthene	22.756	252	20668	0.42	ng/ul	100
26) Benzo(a)pyrene	23.277	252	16805	0.41	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.605	276	23623	0.44	ng/ul	100
28) Dibenzo(a,h)anthracene	25.625	278	18797	0.43	ng/ul	100
29) Benzo(g,h,i)perylene	26.282	276	21128	0.46	ng/ul	100

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

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