

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062422\
 Data File : BN020380.D
 Acq On : 23 Jun 2022 20:24
 Operator : CG/JU
 Sample : SST0.411
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.4211

Quant Time: Jun 24 00:03:26 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 23 23:59:02 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.754	152	3003	0.400	ng/u1	0.00
4) Naphthalene-d8	10.533	136	9785	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.380	164	6571	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.119	188	15259	0.400	ng/u1	0.00
17) Chrysene-d12	21.309	240	13271	0.400	ng/u1	0.00
23) Perylene-d12	23.595	264	10109	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	1416	0.413	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.128	152	6448	0.398	ng/u1	0.00
18) Fluoranthene-d10	19.149	212	16832	0.614	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.239	88	1402	0.425	ng/u1	100
5) Naphthalene	10.583	128	10611	0.339	ng/u1	100
7) 2-Methylnaphthalene	12.205	142	7154	0.353	ng/u1	100
8) 1-Methylnaphthalene	12.420	142	7713	0.387	ng/u1	100
10) Acenaphthylene	14.098	152	10859	0.346	ng/u1	100
11) Acenaphthene	14.440	153	8766	0.292	ng/u1	100
12) Fluorene	15.430	166	10434	0.335	ng/u1	100
14) Pentachlorophenol	16.785	266	1050	0.358	ng/u1	100
15) Phenanthrene	17.161	178	18438	0.334	ng/u1	100
16) Anthracene	17.254	178	15830	0.339	ng/u1	100
19) Fluoranthene	19.181	202	20613	0.385	ng/u1	100
20) Pyrene	19.544	202	21519	0.371	ng/u1	100
21) Benzo(a)anthracene	21.294	228	16370	0.318	ng/u1	100
22) Chrysene	21.347	228	18622	0.290	ng/u1	100
24) Benzo(b)fluoranthene	22.905	252	16512	0.229	ng/u1	100
25) Benzo(k)fluoranthene	22.951	252	16470	0.245	ng/u1	100
26) Benzo(a)pyrene	23.492	252	14885	0.325	ng/u1	100
27) Indeno(1,2,3-cd)pyrene	25.933	276	14799	0.199	ng/u1	100
28) Dibenzo(a,h)anthracene	25.950	278	12005	0.202	ng/u1	100
29) Benzo(g,h,i)perylene	26.647	276	12767	0.194	ng/u1	100

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

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