

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062223\
 Data File : BN025921.D
 Acq On : 21 Jun 2023 20:22
 Operator : MA/JU
 Sample : SST0.208
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.2224

Quant Time: Jun 22 01:20:23 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 22 01:10:16 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.984	152	7917	0.400	ng/ul	0.00
4) Naphthalene-d8	10.798	136	22580	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.621	164	13146	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.368	188	27741	0.400	ng/ul	0.00
17) Chrysene-d12	21.549	240	18584	0.400	ng/ul	0.00
23) Perylene-d12	24.019	264	17370	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.339	96	1807	0.232	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.387	152	6789	0.237	ng/ul	0.00
18) Fluoranthene-d10	19.391	212	14183	0.242	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.377	88	2022	0.243	ng/ul#	87
5) Naphthalene	10.847	128	12838	0.224	ng/ul	99
7) 2-Methylnaphthalene	12.458	142	8077	0.234	ng/ul	100
8) 1-Methylnaphthalene	12.673	142	8447	0.222	ng/ul	99
10) Acenaphthylene	14.343	152	12086	0.233	ng/ul	99
11) Acenaphthene	14.681	153	9997	0.231	ng/ul	98
12) Fluorene	15.671	166	11497	0.238	ng/ul	99
14) Pentachlorophenol	17.022	266	2047	0.686	ng/ul	98
15) Phenanthrene	17.410	178	18290	0.226	ng/ul	99
16) Anthracene	17.503	178	15937	0.233	ng/ul	97
19) Fluoranthene	19.424	202	19748	0.232	ng/ul	99
20) Pyrene	19.786	202	20880	0.237	ng/ul	99
21) Benzo(a)anthracene	21.535	228	15103	0.255	ng/ul	99
22) Chrysene	21.590	228	15874	0.234	ng/ul	99
24) Benzo(b)fluoranthene	23.265	252	14464	0.204	ng/ul	80
25) Benzo(k)fluoranthene	23.315	252	15203	0.213	ng/ul#	78
26) Benzo(a)pyrene	23.914	252	12564	0.209	ng/ul#	73
27) Indeno(1,2,3-cd)pyrene	26.604	276	13081	0.191	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.638	278	9795	0.188	ng/ul#	74
29) Benzo(g,h,i)perylene	27.396	276	11804	0.192	ng/ul#	92

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

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