

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082621\
 Data File : BN016062.D
 Acq On : 25 Aug 2021 17:24
 Operator : CG/JU
 Sample : SST0.403
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.4203

Quant Time: Aug 25 18:14:18 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN082621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 25 18:14:04 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.871	152	1253	0.400	ng/ul	0.00
4) Naphthalene-d8	10.666	136	5145	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.501	164	3052	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.246	188	6751	0.400	ng/ul	0.00
17) Chrysene-d12	21.442	240	6837	0.400	ng/ul	0.00
23) Perylene-d12	23.836	264	5892	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	714	0.506	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	3138	0.389	ng/ul	0.00
18) Fluoranthene-d10	19.280	212	7470	0.332	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.353	88	680	0.446	ng/ul	100
5) Naphthalene	10.716	128	5732	0.349	ng/ul	100
7) 2-Methylnaphthalene	12.333	142	3888	0.362	ng/ul	100
8) 1-Methylnaphthalene	12.547	142	3601	0.347	ng/ul	100
10) Acenaphthylene	14.224	152	4609	0.292	ng/ul	100
11) Acenaphthene	14.566	153	3963	0.328	ng/ul	100
12) Fluorene	15.552	166	4709	0.352	ng/ul	100
14) Pentachlorophenol	16.900	266	611	0.409	ng/ul	100
15) Phenanthrene	17.288	178	7831	0.320	ng/ul	100
16) Anthracene	17.381	178	6829	0.316	ng/ul	100
19) Fluoranthene	19.308	202	9073	0.287	ng/ul	100
20) Pyrene	19.671	202	9204	0.290	ng/ul	100
21) Benzo(a)anthracene	21.424	228	9055	0.358	ng/ul	100
22) Chrysene	21.477	228	9853	0.325	ng/ul	100
24) Benzo(b)fluoranthene	23.105	252	9397	0.341	ng/ul	100
25) Benzo(k)fluoranthene	23.152	252	9611	0.309	ng/ul	100
26) Benzo(a)pyrene	23.728	252	7916	0.342	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	26.307	276	10373	0.371	ng/ul	100
28) Dibenzo(a,h)anthracene	26.327	278	8661	0.396	ng/ul	100
29) Benzo(g,h,i)perylene	27.068	276	9041	0.349	ng/ul	100

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

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