

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042023\
 Data File : BN025022.D
 Acq On : 20 Apr 2023 04:46
 Operator : CG/JU
 Sample : SST0.295
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.2210

Quant Time: Apr 20 07:47:06 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN042023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 20 07:43:17 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.933	152	7237	0.400	ng/ul	0.00
4) Naphthalene-d8	10.742	136	22291	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.575	164	12843	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.326	188	26606	0.400	ng/ul	0.00
17) Chrysene-d12	21.515	240	22511	0.400	ng/ul	0.00
23) Perylene-d12	23.956	264	18207	0.400	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.309	96	1522	0.193	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.332	152	5442	0.190	ng/ul	0.00
18) Fluoranthene-d10	19.350	212	11585	0.207	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.343	88	1626	0.187	ng/ul#	90
5) Naphthalene	10.792	128	11023	0.194	ng/ul	99
7) 2-Methylnaphthalene	12.403	142	6740	0.191	ng/ul	99
8) 1-Methylnaphthalene	12.623	142	7192	0.192	ng/ul	99
10) Acenaphthylene	14.298	152	9148	0.190	ng/ul	99
11) Acenaphthene	14.635	153	7786	0.189	ng/ul	99
12) Fluorene	15.625	166	8542	0.185	ng/ul	98
14) Pentachlorophenol	16.980	266	845	0.149	ng/ul	98
15) Phenanthrene	17.364	178	14632	0.195	ng/ul	99
16) Anthracene	17.461	178	11515	0.185	ng/ul	99
19) Fluoranthene	19.382	202	15773	0.203	ng/ul	98
20) Pyrene	19.745	202	16299	0.205	ng/ul	99
21) Benzo(a)anthracene	21.497	228	12908	0.182	ng/ul	100
22) Chrysene	21.550	228	14311	0.186	ng/ul	100
24) Benzo(b)fluoranthene	23.207	252	14204	0.194	ng/ul	93
25) Benzo(k)fluoranthene	23.254	252	13595	0.191	ng/ul#	93
26) Benzo(a)pyrene	23.853	252	11935	0.197	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	26.511	276	13432	0.188	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.531	278	10385	0.185	ng/ul	95
29) Benzo(g,h,i)perylene	27.289	276	11519	0.191	ng/ul	96

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

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