

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071624\
 Data File : BN032807.D
 Acq On : 18 Jul 2024 08:57
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 80 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4266

Quant Time: Jul 18 09:24:43 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN071224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 17 02:35:25 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.544	152	4543	0.400	ng/ul	0.00
4) Naphthalene-d8	10.316	136	14685	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.193	164	7960	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.961	188	14410	0.400	ng/ul	0.00
17) Chrysene-d12	21.174	240	13196	0.400	ng/ul	0.00
23) Perylene-d12	23.371	264	16306	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.097	96	2138	0.392	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.922	152	8046	0.363	ng/ul	-0.01
18) Fluoranthene-d10	18.999	212	14029	0.326	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.131	88	2228	0.359	ng/ul#	86
5) Naphthalene	10.365	128	14641	0.371	ng/ul	99
7) 2-Methylnaphthalene	11.993	142	8854	0.360	ng/ul	98
8) 1-Methylnaphthalene	12.213	142	9402	0.369	ng/ul	99
10) Acenaphthylene	13.915	152	13057	0.406	ng/ul	99
11) Acenaphthene	14.258	153	9610	0.361	ng/ul	99
12) Fluorene	15.262	166	9835	0.324	ng/ul	99
14) Pentachlorophenol	16.645	266	1045	0.413	ng/ul	99
15) Phenanthrene	17.003	178	13781	0.351	ng/ul	100
16) Anthracene	17.100	178	12807	0.399	ng/ul	99
19) Fluoranthene	19.031	202	16994	0.317	ng/ul	99
20) Pyrene	19.399	202	17757	0.320	ng/ul	99
21) Benzo(a)anthracene	21.159	228	16425	0.391	ng/ul	98
22) Chrysene	21.212	228	19119	0.315	ng/ul	98
24) Benzo(b)fluoranthene	22.711	252	18642	0.324	ng/ul	94
25) Benzo(k)fluoranthene	22.754	252	20328	0.290	ng/ul	95
26) Benzo(a)pyrene	23.275	252	18651	0.375	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.586	276	23383	0.347	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.586	278	18908	0.362	ng/ul#	95
29) Benzo(g,h,i)perylene	26.267	276	19182	0.340	ng/ul	97

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

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