

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102221\
 Data File : BN017037.D
 Acq On : 21 Oct 2021 08:46
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4274

Quant Time: Oct 21 10:29:49 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN101921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 21 10:28:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.544	152	2259	0.400	ng/ul	0.00
4) Naphthalene-d8	10.317	136	8272	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.199	164	4632	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.953	188	10573	0.400	ng/ul	0.00
17) Chrysene-d12	21.165	240	11646	0.400	ng/ul	0.00
23) Perylene-d12	23.375	264	12935	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.076	96	965	0.393	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.923	152	4305	0.365	ng/ul	0.00
18) Fluoranthene-d10	18.990	212	10693	0.372	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.110	88	941	0.318	ng/ul#	48
5) Naphthalene	10.366	128	8816	0.352	ng/ul	100
7) 2-Methylnaphthalene	11.994	142	5717	0.367	ng/ul	100
8) 1-Methylnaphthalene	12.214	142	5881	0.370	ng/ul	100
10) Acenaphthylene	13.912	152	7294	0.390	ng/ul	99
11) Acenaphthene	14.259	153	6342	0.395	ng/ul	98
12) Fluorene	15.253	166	7137	0.384	ng/ul	100
14) Pentachlorophenol	16.611	266	727	0.285	ng/ul	99
15) Phenanthrene	16.991	178	13008	0.387	ng/ul	100
16) Anthracene	17.084	178	10402	0.367	ng/ul	99
19) Fluoranthene	19.023	202	15355	0.385	ng/ul	98
20) Pyrene	19.385	202	15481	0.377	ng/ul	100
21) Benzo(a)anthracene	21.148	228	13565	0.357	ng/ul	100
22) Chrysene	21.200	228	17466	0.404	ng/ul	99
24) Benzo(b)fluoranthene	22.711	252	18076	0.371	ng/ul	99
25) Benzo(k)fluoranthene	22.755	252	21331	0.398	ng/ul	98
26) Benzo(a)pyrene	23.278	252	15824	0.364	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.606	276	24832	0.402	ng/ul	99
28) Dibenzo(a,h)anthracene	25.627	278	19591	0.399	ng/ul	99
29) Benzo(g,h,i)perylene	26.281	276	21566	0.403	ng/ul	100

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

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