

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102221\
 Data File : BN017070.D
 Acq On : 22 Oct 2021 13:15
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4277

Quant Time: Oct 22 13:46:36 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 : Fri Oct 22 10:13:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.548	152	592	0.400	ng/ul	0.00
4) Naphthalene-d8	10.321	136	2148	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.197	164	1243	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.952	188	3021	0.400	ng/ul	0.00
17) Chrysene-d12	21.164	240	5752	0.400	ng/ul	0.00
23) Perylene-d12	23.376	264	9257	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.076	96	245	0.381	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.921	152	1085	0.354	ng/ul	0.00
18) Fluoranthene-d10	18.993	212	3896	0.274	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.109	88	215	0.278	ng/ul#	88
5) Naphthalene	10.370	128	2327	0.358	ng/ul#	94
7) 2-Methylnaphthalene	11.998	142	1461	0.361	ng/ul	100
8) 1-Methylnaphthalene	12.218	142	1556	0.377	ng/ul	100
10) Acenaphthylene	13.915	152	2197	0.437	ng/ul	96
11) Acenaphthene	14.262	153	1705	0.395	ng/ul	97
12) Fluorene	15.257	166	1941	0.389	ng/ul#	99
14) Pentachlorophenol	16.610	266	326	0.448	ng/ul	99
15) Phenanthrene	16.994	178	3748	0.391	ng/ul	98
16) Anthracene	17.087	178	3210	0.397	ng/ul	94
19) Fluoranthene	19.021	202	5481	0.278	ng/ul	98
20) Pyrene	19.388	202	6058	0.299	ng/ul	100
21) Benzo(a)anthracene	21.149	228	7472	0.398	ng/ul	99
22) Chrysene	21.202	228	8547	0.401	ng/ul	99
24) Benzo(b)fluoranthene	22.712	252	12158	0.349	ng/ul	99
25) Benzo(k)fluoranthene	22.756	252	13360	0.348	ng/ul	97
26) Benzo(a)pyrene	23.280	252	12059	0.387	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.608	276	20498	0.463	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.625	278	15841	0.451	ng/ul	99
29) Benzo(g,h,i)perylene	26.282	276	17526	0.457	ng/ul	100

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

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