

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052022\
 Data File : BN019891.D
 Acq On : 21 May 2022 05:01
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4360

Quant Time: May 26 05:15:59 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN042922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 17 01:41:05 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	3355	0.400	ng/ul	0.00
4) Naphthalene-d8	10.639	136	11721	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.474	164	7535	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.213	188	16074	0.400	ng/ul #	0.00
17) Chrysene-d12	21.397	240	15035	0.400	ng/ul	0.00
23) Perylene-d12	23.735	264	12403	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	1509	0.380	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.229	152	8458	0.477	ng/ul	0.00
18) Fluoranthene-d10	19.239	212	19213	0.404	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.353	88	1605	0.381	ng/ul	92
5) Naphthalene	10.689	128	14745	0.389	ng/ul	99
7) 2-Methylnaphthalene	12.300	142	10054	0.435	ng/ul#	100
8) 1-Methylnaphthalene	12.520	142	10081	0.471	ng/ul#	100
10) Acenaphthylene	14.192	152	16053	0.431	ng/ul	99
11) Acenaphthene	14.534	153	11391	0.380	ng/ul	99
12) Fluorene	15.520	166	13102	0.399	ng/ul	99
14) Pentachlorophenol	16.867	266	1280	0.370	ng/ul	90
15) Phenanthrene	17.251	178	21934	0.371	ng/ul	99
16) Anthracene	17.344	178	20195	0.419	ng/ul	98
19) Fluoranthene	19.267	202	25112	0.365	ng/ul	99
20) Pyrene	19.629	202	25701	0.374	ng/ul	100
21) Benzo(a)anthracene	21.382	228	22506	0.401	ng/ul	99
22) Chrysene	21.435	228	23295	0.352	ng/ul	99
24) Benzo(b)fluoranthene	23.025	252	20887	0.332	ng/ul	92
25) Benzo(k)fluoranthene	23.072	252	20906	0.345	ng/ul	97
26) Benzo(a)pyrene	23.630	252	19613	0.366	ng/ul#	78
27) Indeno(1,2,3-cd)pyrene	26.138	276	19766	0.296	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.154	278	16288	0.291	ng/ul#	80
29) Benzo(g,h,i)perylene	26.872	276	16345	0.267	ng/ul	97

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

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