

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110923\
 Data File : BN028574.D
 Acq On : 11 Nov 2023 01:45
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4308

Quant Time: Nov 11 02:33:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 09 16:30:32 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.703	152	9300	0.400	ng/ul	0.00
4) Naphthalene-d8	10.491	136	31017	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.349	164	19657	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.099	188	45840	0.400	ng/ul	0.00
17) Chrysene-d12	21.311	240	36589	0.400	ng/ul	0.00
23) Perylene-d12	23.611	264	34943	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.151	96	4180	0.407	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.091	152	19791	0.424	ng/ul	0.00
18) Fluoranthene-d10	19.137	212	51513	0.437	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.184	88	4305	0.379	ng/ul#	88
5) Naphthalene	10.540	128	35553	0.409	ng/ul	100
7) 2-Methylnaphthalene	12.162	142	24130	0.421	ng/ul	99
8) 1-Methylnaphthalene	12.382	142	24612	0.423	ng/ul	100
10) Acenaphthylene	14.067	152	40849	0.416	ng/ul	100
11) Acenaphthene	14.409	153	29967	0.415	ng/ul	99
12) Fluorene	15.404	166	34651	0.413	ng/ul	99
14) Pentachlorophenol	16.752	266	4204	0.382	ng/ul	99
15) Phenanthrene	17.141	178	57781	0.410	ng/ul	100
16) Anthracene	17.234	178	52517	0.407	ng/ul	100
19) Fluoranthene	19.169	202	67424	0.433	ng/ul	97
20) Pyrene	19.532	202	69752	0.432	ng/ul	100
21) Benzo(a)anthracene	21.293	228	58382	0.404	ng/ul	99
22) Chrysene	21.346	228	57358	0.406	ng/ul	100
24) Benzo(b)fluoranthene	22.915	252	53959	0.411	ng/ul	100
25) Benzo(k)fluoranthene	22.962	252	57531	0.419	ng/ul	98
26) Benzo(a)pyrene	23.509	252	48577	0.392	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.971	276	56075	0.408	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.995	278	44977	0.410	ng/ul	96
29) Benzo(g,h,i)perylene	26.696	276	45506	0.404	ng/ul	99

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

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