

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032524\
 Data File : BN030496.D
 Acq On : 26 Mar 2024 00:47
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4364

Quant Time: Mar 26 01:14:34 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN032524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 25 13:38:40 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.724	152	8932	0.400	ng/u1	0.00
4) Naphthalene-d8	10.507	136	25599	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.359	164	14027	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.109	188	31707	0.400	ng/u1	0.00
17) Chrysene-d12	21.299	240	22841	0.400	ng/u1	0.00
23) Perylene-d12	23.546	264	20744	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.218	96	4319	0.412	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.102	152	14812	0.411	ng/u1	0.00
18) Fluoranthene-d10	19.138	212	30631	0.430	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	4484	0.413	ng/u1	94
5) Naphthalene	10.557	128	29788	0.408	ng/u1	99
7) 2-Methylnaphthalene	12.173	142	19464	0.411	ng/u1	100
8) 1-Methylnaphthalene	12.393	142	20615	0.412	ng/u1	100
10) Acenaphthylene	14.082	152	28759	0.404	ng/u1	99
11) Acenaphthene	14.424	153	22257	0.406	ng/u1	98
12) Fluorene	15.414	166	25572	0.408	ng/u1	98
14) Pentachlorophenol	16.750	266	2319	0.389	ng/u1	99
15) Phenanthrene	17.147	178	43016	0.404	ng/u1	99
16) Anthracene	17.239	178	36208	0.403	ng/u1	100
19) Fluoranthene	19.171	202	46389	0.427	ng/u1	99
20) Pyrene	19.533	202	45455	0.424	ng/u1	100
21) Benzo(a)anthracene	21.284	228	37117	0.394	ng/u1	99
22) Chrysene	21.334	228	40768	0.409	ng/u1	99
24) Benzo(b)fluoranthene	22.871	252	37111	0.378	ng/u1	98
25) Benzo(k)fluoranthene	22.915	252	40916	0.408	ng/u1	98
26) Benzo(a)pyrene	23.450	252	30390	0.387	ng/u1	99
27) Indeno(1,2,3-cd)pyrene	25.830	276	38011	0.399	ng/u1	100
28) Dibenzo(a,h)anthracene	25.847	278	30443	0.400	ng/u1	98
29) Benzo(g,h,i)perylene	26.521	276	29926	0.363	ng/u1	99

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

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