

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN103122\
 Data File : BN022408.D
 Acq On : 31 Oct 2022 20:00
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4332

Quant Time: Nov 01 01:40:39 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 01 01:09:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.932	152	3334	0.400	ng/u1	0.00
4) Naphthalene-d8	10.756	136	10807	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.600	164	6051	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.357	188	12018	0.400	ng/u1	0.00
17) Chrysene-d12	21.558	240	7385	0.400	ng/u1	0.00
23) Perylene-d12	24.005	264	5139	0.400	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	1686	0.389	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.351	152	6348	0.388	ng/u1	0.00
18) Fluoranthene-d10	19.393	212	10748	0.434	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.278	88	1808	0.406	ng/u1	92
5) Naphthalene	10.805	128	12477	0.391	ng/u1	100
7) 2-Methylnaphthalene	12.428	142	7763	0.386	ng/u1	100
8) 1-Methylnaphthalene	12.648	142	8010	0.387	ng/u1	100
10) Acenaphthylene	14.322	152	10329	0.380	ng/u1	99
11) Acenaphthene	14.665	153	8800	0.384	ng/u1	99
12) Fluorene	15.655	166	9640	0.362	ng/u1	99
14) Pentachlorophenol	17.011	266	1182	0.414	ng/u1	100
15) Phenanthrene	17.400	178	15079	0.374	ng/u1	99
16) Anthracene	17.493	178	12420	0.365	ng/u1	99
19) Fluoranthene	19.421	202	14580	0.431	ng/u1	100
20) Pyrene	19.788	202	14478	0.426	ng/u1	99
21) Benzo(a)anthracene	21.541	228	10385	0.375	ng/u1	99
22) Chrysene	21.596	228	11091	0.378	ng/u1	99
24) Benzo(b)fluoranthene	23.254	252	9982	0.387	ng/u1	91
25) Benzo(k)fluoranthene	23.303	252	9405	0.375	ng/u1#	94
26) Benzo(a)pyrene	23.897	252	8216	0.399	ng/u1#	84
27) Indeno(1,2,3-cd)pyrene	26.571	276	10163	0.397	ng/u1	97
28) Dibenzo(a,h)anthracene	26.595	278	7940	0.389	ng/u1	95
29) Benzo(g,h,i)perylene	27.363	276	8354	0.401	ng/u1	98

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

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