

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032223\
 Data File : BN024500.D
 Acq On : 23 Mar 2023 04:15
 Operator : CG/JU
 Sample : PB151557BS
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS557

Quant Time: Mar 23 04:44:35 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN030923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 23 03:32:22 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.025	152	9285	0.400	ng/u1	0.00
4) Naphthalene-d8	10.839	136	26865	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.661	164	13869	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.409	188	29629	0.400	ng/u1	0.00
17) Chrysene-d12	21.585	240	22701	0.400	ng/u1	0.00
23) Perylene-d12	24.082	264	20373	0.400	ng/u1	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	7056	0.691	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.423	152	13673	0.428	ng/u1	0.00
18) Fluoranthene-d10	19.431	212	27693	0.458	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.427	88	18096	1.628	ng/u1	90
5) Naphthalene	10.889	128	28019	0.418	ng/u1	99
7) 2-Methylnaphthalene	12.495	142	17191	0.423	ng/u1	99
8) 1-Methylnaphthalene	12.715	142	18033	0.422	ng/u1	100
10) Acenaphthylene	14.384	152	24576	0.425	ng/u1	100
11) Acenaphthene	14.726	153	18599	0.417	ng/u1	98
12) Fluorene	15.707	166	20608	0.414	ng/u1	97
14) Pentachlorophenol	17.058	266	5323	0.647	ng/u1#	72
15) Phenanthrene	17.447	178	34122	0.402	ng/u1	100
16) Anthracene	17.540	178	30119	0.404	ng/u1	100
19) Fluoranthene	19.459	202	39323	0.435	ng/u1	100
20) Pyrene	19.822	202	40676	0.444	ng/u1	99
21) Benzo(a)anthracene	21.571	228	32870	0.444	ng/u1	100
22) Chrysene	21.626	228	34205	0.437	ng/u1	99
24) Benzo(b)fluoranthene	23.313	252	33449	0.430	ng/u1	96
25) Benzo(k)fluoranthene	23.369	252	31684	0.392	ng/u1	95
26) Benzo(a)pyrene	23.968	252	27388	0.409	ng/u1	93
27) Indeno(1,2,3-cd)pyrene	26.693	276	31296	0.415	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.716	278	23769	0.415	ng/u1	96
29) Benzo(g,h,i)perylene	27.501	276	27426	0.419	ng/u1	97

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

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