

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032223\
 Data File : BN024498.D
 Acq On : 23 Mar 2023 03:01
 Operator : CG/JU
 Sample : PB151553BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS553

Quant Time: Mar 23 03:34:39 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	10776	0.400	ng/u1	0.00
4) Naphthalene-d8	10.839	136	31165	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.661	164	16309	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.409	188	34965	0.400	ng/u1	0.00
17) Chrysene-d12	21.586	240	27010	0.400	ng/u1	0.00
23) Perylene-d12	24.082	264	24483	0.400	ng/u1	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	7222	0.609	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.423	152	14272	0.385	ng/u1	0.00
18) Fluoranthene-d10	19.432	212	29704	0.413	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.427	88	18678	1.448	ng/u1	91
5) Naphthalene	10.889	128	29011	0.373	ng/u1	100
7) 2-Methylnaphthalene	12.495	142	17894	0.379	ng/u1	100
8) 1-Methylnaphthalene	12.715	142	18851	0.380	ng/u1	100
10) Acenaphthylene	14.384	152	26055	0.383	ng/u1	99
11) Acenaphthene	14.721	153	19499	0.372	ng/u1	98
12) Fluorene	15.707	166	21719	0.371	ng/u1	98
14) Pentachlorophenol	17.059	266	6039	0.622	ng/u1#	72
15) Phenanthrene	17.447	178	35719	0.357	ng/u1	100
16) Anthracene	17.540	178	32163	0.366	ng/u1	99
19) Fluoranthene	19.459	202	41597	0.387	ng/u1	100
20) Pyrene	19.822	202	43293	0.397	ng/u1	99
21) Benzo(a)anthracene	21.568	228	35773	0.406	ng/u1	100
22) Chrysene	21.627	228	36238	0.389	ng/u1	99
24) Benzo(b)fluoranthene	23.316	252	34685	0.371	ng/u1	95
25) Benzo(k)fluoranthene	23.372	252	33952	0.350	ng/u1	94
26) Benzo(a)pyrene	23.968	252	28830	0.358	ng/u1	93
27) Indeno(1,2,3-cd)pyrene	26.696	276	32300	0.356	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.720	278	24916	0.362	ng/u1	97
29) Benzo(g,h,i)perylene	27.505	276	27853	0.354	ng/u1	97

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

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