

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062422\
 Data File : BN020399.D
 Acq On : 24 Jun 2022 16:01
 Operator : CG/JU
 Sample : PB145637BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS637

Quant Time: Jun 25 00:55:05 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 00:17:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.750	152	2553	0.400	ng/u1	0.00
4) Naphthalene-d8	10.535	136	8591	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.382	164	5768	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.120	188	13357	0.400	ng/u1	0.00
17) Chrysene-d12	21.309	240	11945	0.400	ng/u1	0.00
23) Perylene-d12	23.592	264	9640	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	2167	0.733	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.130	152	5421	0.385	ng/u1	0.00
18) Fluoranthene-d10	19.151	212	15857	0.403	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.239	88	7055	2.336	ng/u1#	82
5) Naphthalene	10.584	128	11921	0.450	ng/u1	99
7) 2-Methylnaphthalene	12.201	142	7997	0.450	ng/u1	99
8) 1-Methylnaphthalene	12.421	142	8093	0.481	ng/u1	100
10) Acenaphthylene	14.099	152	12402	0.454	ng/u1	99
11) Acenaphthene	14.442	153	9683	0.442	ng/u1	99
12) Fluorene	15.427	166	11551	0.445	ng/u1	100
14) Pentachlorophenol	16.782	266	2025	0.752	ng/u1	100
15) Phenanthrene	17.162	178	20696	0.446	ng/u1	100
16) Anthracene	17.251	178	18249	0.455	ng/u1	99
19) Fluoranthene	19.179	202	24511	0.449	ng/u1	100
20) Pyrene	19.541	202	25505	0.451	ng/u1	98
21) Benzo(a)anthracene	21.295	228	21737	0.507	ng/u1	99
22) Chrysene	21.347	228	23935	0.501	ng/u1	100
24) Benzo(b)fluoranthene	22.905	252	22132	0.496	ng/u1	98
25) Benzo(k)fluoranthene	22.946	252	22412	0.498	ng/u1	98
26) Benzo(a)pyrene	23.492	252	20497	0.506	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	25.926	276	20600	0.511	ng/u1#	98
28) Dibenzo(a,h)anthracene	25.940	278	16769	0.515	ng/u1	99
29) Benzo(g,h,i)perylene	26.641	276	17532	0.506	ng/u1	99

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

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